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TECHNICAL ASSISTANCE TEAM FOR EMERGENCY RESPONSE REMOVAL AND PREVENTION
EPA CONTRACT 68-WO-0036

TO: Dave Turner, Remedial Project Manager, EPA
Philadelphia, Pennsylvania

FROM: John J. Mueck, Jr., Technical Assistance Team, Region
III, Wheeling, West Virginia *CAH*

THRU: Joseph B. Carter, ATATL, Region III, *JC*
Wheeling, West Virginia

THRU: Gerald T. Heston, OSC, EPA Region III
Philadelphia, Pennsylvania

SUBJECT: Trip Report - Westinghouse Sharon NPL Site,
Sharon, Mercer County, Pennsylvania
TDD# 9410-160 PCS# 1160

DATE: November 5, 1994

Background

The Sharon Westinghouse NPL Site is located at 369 Sharpsville Avenue in Sharon, Mercer County, Pennsylvania. For 63 years, Westinghouse Electric Corporation owned and operated this Site where they manufactured, repaired, and shipped electrical transformers and dielectric components. Westinghouse used polychlorinated biphenyls (PCBs) as a dielectric fluid in the manufactured components that they produced at the Site. A number of chemicals were used in the manufacturing process, and various waste streams were generated containing PCBs, metals, and solvents. During operations, materials, including PCB-contaminated oils, were routinely burned in an incinerator located onsite. In 1985, Westinghouse shut down operations and sold some portions of their property to neighboring industries. The Site was proposed for the National Priorities List (NPL) in 1988.

Currently, Westinghouse is performing a Remedial Investigation/Feasibility Study (RI/FS) at the Site under an Order with the Commonwealth of Pennsylvania. Samples taken during the RI have shown that contaminated soil is present on the Site. EPA determined that off-site soil sampling was necessary to investigate migration of these contaminants into the residential, commercial and industrial properties surrounding the Site.

by F. Weston, Inc.

MAJOR PROGRAMS DIVISION

In Association with Foster Wheeler Enviresponse, Inc., Resource Applications, Inc., C.C. Johnson & Malhotra, P.C., R.E. Sarriera Associates, and GRB Environmental Services, Inc.

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In May 1994, U.S. EPA On-Scene Coordinator (OSC) Heston tasked Roy F. Weston, Inc., Technical Assistance Team (TAT) to collect 31 off-site soil samples for analysis for PCBs, pesticides, metals, cyanide, volatile organics, semi-volatile organics, dioxins, and dibenzo furans.

Sampling Activities

On August 2, 1994, four members of TAT mobilized from the Wheeling, West Virginia office to the Westinghouse Sharon Site to conduct sampling activities. TAT personnel met with EPA Remedial Project Manager (RPM) for the Site, Dave Turner, and EPA Office of Public Affairs (OPA), Pat Gaughan. RPM Turner and OPA Gaughan had prepared access agreements for property owners which allowed EPA to access and sample individual yards. Table 1 in Attachment B of this report lists the sample locations, residents and addresses where access was granted.

During the first day of the sampling event, August 2, 1994, TAT personnel signed in at the guard shack at the Westinghouse property entrance off Sharpesville Avenue and set up a support zone and a sample preparation area inside the Westinghouse Facility gate. TAT used their Emergency Response Vehicle (ERV) as a base of operations and to supply power to the blenders used in preparing dioxin samples. In addition, the office space was used to complete sampling paperwork.

TAT personnel separated into two teams, a sampling team and a sample preparation team, and began sampling activities. A trip blank, a blender cup rinsate blank, and soil samples S01 - S12 were obtained throughout the day. In addition, sample S06 was a double volume sample for laboratory quality assurance (QA). The soil samples were obtained from 3" - 6" depth following the guidelines of the sampling plan with one exception. Stainless steel scoops were used for sample collection instead of dedicated plastic scoops. The stainless steel scoops were decontaminated after each use. Samples were obtained from locations showing minimal recent disturbance and minimal potential for cross-contamination from other sources (i.e. vehicles). Where available, residents assisted TAT by showing locations of their properties which had been undisturbed. The sampling team marked the locations of the samples on a field map, measured and logged distances from the sampling point to nearby landmarks, and photographed the sampling locations. The sampling locations are shown on the sample location map in Attachment A of this report.

The sample preparation team followed the guidelines of the sampling plan in preparing the dioxin/dibenzo furan samples with one exception. The samples were first obtained in clean sample jars

then transferred to the stainless steel blending cup for preparation. After blending, the soil sample was returned to the original container. Five soil samples (S01, S04, S06, S07, S09) and one blender cup rinsate blank were prepared for dioxin/dibenzo furan analysis by the end of the day.

On the second day of the sampling event, August 3, 1994, TAT personnel continued and completed sampling and sample preparation activities. Two stainless steel spoon rinsate samples (RB-1, RB-2) and soil samples S13 - S31 were obtained. Of the soil samples collected, ten were comprised of additional volumes for dioxin/dibenzo furan analysis (S15, S16, S17, S20, S24, S27, S28, S29, S30, S31), four were from background locations (S28, S29, S30, S31), sample S16 was a blind duplicate of S15, and sample S23 was a blind duplicate of sample S22.

TAT personnel departed the Site on August 3, 1994 and returned to the Wheeling, WV office on the morning of August 4, 1994. Contract Laboratory Program (CLP) samples were shipped to the CLP designated labs on August 4, 1994. The dioxin/dibenzo furan samples were shipped on August 5, 1994. TAT received the analytical data by October 1994.

Sample Analysis

As shown in Table 1, one trip blank, two scoop rinsate samples, one blender cup rinsate, and 31 soil samples (including two blind duplicates and two extra volume samples) were obtained during the trip. All samples except the trip blank and blender cup rinsate blank were analyzed for total metals, cyanide, volatile organics (VOA), base, neutral and acid extractables (BNA), and pesticides/polychlorinated biphenyls (PEST/PCB) through the Contract Laboratory Program (CLP). The trip blank was analyzed for VOA. Through the CLP, Mack Laboratories, Inc., Pittsburgh, PA performed the total metals and cyanide analyses (see Attachment C). EnviroSystems, Inc., Columbia, MD performed the VOA, BNA and PEST/PCB analyses (see Attachment D).

Fifteen of the 31 soil samples and the blender cup rinsate sample were analyzed for isomer specific polychlorinated dibenzo dioxins and polychlorinated dibenzo furans and to determine the toxicity equivalents as 2,3,7,8 tetra chloro dibenzo dioxin (2,3,7,8 TCDD). These analyses were conducted by the Weston Lionville Analytical Laboratory, Lionville, PA (see Attachment E).

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Westinghouse Sharon NPL Site
Trip Report
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Attachments

Attachment A - Sample Location Maps

Attachment B - Sample Information Log

Attachment C - Inorganics Data Validation, Analytical Summary,
Traffic Report

Attachment D - Organics Data Validation, Analytical Summary,
Traffic Report

Attachment E - Dioxin/Dibenzo furan Data Validation, Analytical
Summary, Chain-of-Custody

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Attachment A
Sample Location Maps

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Attachment B
Sample Information Log

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TABLE 1

WESTINGHOUSE - SHARON PA. SAMPLE INFORMATION LOG

EPA/TAT

SAMPLE #	ADDRESS	LOCATION	ANALYSIS					
			DECON CONTAMINANT	TOTAL METAL	Co	VOA	BMA	PCB PEST
TB1	Tip Blank 6/2/94	N/A						
RB1	Rimsale, Sampling Scoops	Sample preparation area						
RB2	Rimsale, Sampling Scoops	Sample preparation area						
	Rimsale, Bleeder Cup	Sample preparation area						
S01		44' south of SE corner of garage, under hedge						
S02		6' west of NE corner of garage, 2' to the N						
S03		3' south of SE corner of porch						
S04		16' east of southern porch steps						
S05		20' west of SW corner of house						
S06/det. vol.		11' south of SE corner of house						
S07		2' north of north end of house						
S08		4' west of SW corner of house						
S09		SE corner of house						
S10		12' north of NE corner of house						
S11		13' west of NW corner of house						
S12		20' west of SW corner of house, along property line						
S13		4' east of center of E side of house						
S14		13' east of NE corner of garage						
S15		3' west of SW corner of porch						
S16		Blind duplicate of S13						
S17		10' east of SE corner of porch						
S18		20' east of SE corner of house						
S19/det. vol.		13' north SE of SW corner of house						
S20		3' south of center garage SE corner at base of tree						
S21		10' east and 20' north of SE corner of building						
S22		30' east of NE corner of house						
S23		Blind duplicate of S22						
S24	Yarnac, Pagan State Campus	3' west of second large tree west of Shamago on Yarnac						
S25	Reno Athletic Field	10' east of westernmost maple on Reno St.						
S26	Alexander Tennis Courts	Northern field, 37' north of center gate						
S27		40' north and 20' east of fire hydrant in western field						
S28		10' east of southern water meter cover in front yard						
S29		23' north of second long window from the west in the rear						
S30		10' west of road, 20' north of driveway						
S31	Same owner, 2 mi. north on Ridge Rd.	430' west and 130' south of field entry gate						

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Attachment C

Inorganics Data Validation, Analytical Summary, Traffic Report

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Environmental Systems & Technologies Co.

Environmental Services Assistance Teams
Region 3
1419 Forest Drive, Suite 104
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Phone: (410) 268-7705
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DATE: SEPTEMBER 23, 1994

SUBJECT: INORGANIC DATA VALIDATION CASE 22533
SITE: WESTINGHOUSE SHARON

FROM: *for* SHOBHA BODDU *HP* MAHBOOBH MECANIC *117*
INORGANIC DATA REVIEWER SENIOR OVERSIGHT CHEMIST

TO: SUSANNE STEVENS
ESAT ACTING REGIONAL PROJECT OFFICER

THROUGH: DALE S. BOSHART *DB*
ESAT TEAM MANAGER

OVERVIEW

The set of samples for Case 22533 consisted of thirty-three (33) soil samples and two (2) associated rinsate blanks. Included in the sample set were two (2) field duplicate pairs. The samples were analyzed by Mack Laboratories (MACK) for total metals and cyanide (CN) according to the Contract Laboratory Program (CLP) Routine Analytical Services (RAS) Statement of Work (SOW) ILM03.0.

The Chemical Health Advisory Level was exceeded for the cadmium (Cd) and lead (Pb) analytes in several samples. The Regional Project Manager (RPM) was notified. See Table 4 for details.

SUMMARY

The samples were analyzed under two (2) Sample Delivery Groups (SDGs) and the data validation was also performed on an SDG basis.

All analytes except antimony (Sb) and selenium (Se) in SDG MCRS73 were successfully analyzed in all samples. Areas of concern with respect to data usability are listed below according to the seriousness of the problem.

MAJOR PROBLEM

The Contract Required detection Limit (CRDL) standard recoveries fall below 50% for the Sb and Se analytes in SDG MCRS73. The quantitation limits in the affected samples are unusable and have been qualified "R".

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MINOR ISSUES

The rinsate (RB), preparation (PB) and continuing calibration (CCB) blanks had reported results greater than the Instrument Detection Limit (IDL) for the analytes listed below. The reported results which are less than five times ($< 5x$) the blank concentrations may be biased high and have been qualified "B".

<u>SDG</u>	<u>Blank Type</u>	<u>Analyte(s)</u>
MCRS57	RB	cadmium (Cd), calcium (Ca), mercury (Hg), nickel (Ni), sodium (Na), zinc (Zn)
	PB (Soil)	potassium (K)
	PB (Aqueous)	Ca, Hg, Na
	CCB	copper (Cu), Ni
MCRS73	RB	Cd, Ca, Cu, Pb, Hg, Ni, Zn
	PB	Na

The matrix spike recovery was high and the CRDL standard recovery was extremely low for the Se analyte in SDG MCRS73, resulting in opposing bias effects. The reported results which are $< 2x$ CRDL for this analyte are estimated and have been qualified "J".

The matrix spike recoveries were low for the Sb and arsenic (As) analytes for the soil samples in SDG MCRS57 and for the Sb analyte in SDG MCRS73. The quantitation limits and reported results for these analytes may be biased low and have respectively been qualified "UL" and "L", unless superseded by the "R" qualifier.

The soil matrix spike recovery was high for the Hg and Se analytes in SDG MCRS57. The reported results for these analytes in the affected samples may be biased high and have been qualified "K", unless superseded by the "B" qualifier.

The CRDL standard recoveries were high for the analytes given below. The reported results which are $< 2x$ CRDL may be biased high and have been qualified "K", unless superseded by the "B" qualifier.

<u>SDG</u>	<u>Analytes</u>
MCRS57	Cd, Cu, Ni, Se
MCRS73	Cd, Ni

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The CRDL standard recoveries were low for the analytes given below. The reported results which are $< 2 \times$ CRDL and quantitation limits for these analytes may be biased low and have been qualified "L" and "UL", respectively.

<u>SDG</u>	<u>Analytes</u>
MCRS57	Sb, As, chromium (Cr)
MCRS73	Cu

The analytical spike recovery was low for the thallium (Tl) analyte in sample MCRS70 of SDG MCRS57. The quantitation limit for this analyte in this sample may be biased low and has been qualified "UL".

The continuing calibration blanks had negative results with absolute values greater than the IDL for the aluminum (Al) and K analytes in SDG MCRS57, and Tl analyte in SDG MCRS73. The quantitation limits for these analytes which have negative raw data results may be biased low and have been qualified "UL".

NOTES

The laboratory marked the sample results for the Pb, Hg and CN analytes in SDG MCRS57, and the CN analyte in SDG MCRS73 with a "*" denoting laboratory duplicate results outside the contractual control limits ($20\% \text{ RPD}, \pm \text{CRDL}$). Since the technical control limits for the soil samples ($35\% \text{ RPD}, \pm 2 \times \text{CRDL}$) were not exceeded for these analytes, no data were qualified during validation.

The laboratory did not account for the 10-fold dilution while reporting the result for the Fe analyte on the Form I for sample MCRS67 (SDG MCRS57). This discrepancy has been rectified during validation.

The results for the field duplicate pairs were within the technical control limits except for the analytes listed in Table 5. Because there are no criteria established in Region III for field duplicate precision, no data were qualified.

The analytical spike recoveries were high for the Pb and Tl analytes in several samples in SDG MCRS57 and for the Se and Tl analytes in SDG MCRS73. These analytes were not detected above the IDL in these samples. Because high recoveries do not affect quantitation limits, no action was taken.

Soil sample results are calculated on the basis of the raw data values (in $\mu\text{g/L}$), the gram weight of sample used, the volume of the digestate, and the % solids according to the following equation:

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$$\text{mg/Kg} = \frac{(\text{raw value, ug/L}) (\text{digestate volume, L})}{(\text{weight, g}) (\% \text{ solids}/100)}$$

To obtain quantitation limit, insert the IDL (Form X) for the raw value; refer to Form XIII to obtain each sample preparation weight and volume used. The quantitation limits thus obtained are specific for each sample and preparation method.

The data were reviewed in accordance with the National Functional Guidelines for Evaluating Inorganic Analyses (IM2 Level), with modifications for use within Region III.

INFORMATION REGARDING REPORT CONTENT

Table 1A is a summary of qualifiers added to the results of the laboratory during validation.

ATTACHMENTS

TABLE 1A	SUMMARY OF QUALIFIERS ON DATA SUMMARY AFTER DATA VALIDATION
TABLE 1B	CODES USED IN COMMENTS COLUMN
TABLE 2	GLOSSARY OF DATA QUALIFIER CODES
TABLE 3	DATA SUMMARY FORM
TABLE 4	SAMPLES EXCEEDING THE CHEMICAL HEALTH ADVISORY LIMITS
TABLE 5	SUMMARY OF FIELD DUPLICATE RESULTS
APPENDIX A	RESULTS REPORTED BY LABORATORY FORM 1s
APPENDIX B	SUPPORT DOCUMENTATION.

DCN: SB409A03.WHS

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TABLE 1A
SUMMARY OF QUALIFIERS ON DATA SUMMARY
FORM AFTER DATA VALIDATION

<u>ANALYTE</u>	<u>SDG</u>	<u>SAMPLES AFFECTED+</u>	<u>POSITIVE VALUES</u>	<u>NON- DETECTED VALUES</u>	<u>BIAS</u>	<u>COMMENTS*</u>
Al	MCRS57	70,81		UL	Low	CCN(-167 µg/L)
Sb	MCRS57	70,81		UL	Low	CRL(79.2%,53.3%)
		57-60		UL	Low	CRL(79.2%,53.3%) MSL(70.6%)
		61-69,71,72	L	UL	Low	CRL(81.7%) MSL(70.6%)
	MCRS73	All samples		R	Extr. Low	CRE(43.3%) MSL(43.0%)
As	MCRS57	70,81		UL	Low	CRL(77.0%)
		All soil samples	L		Low	MSL(64.5%)
Cd	MCRS57	All soil samples except 57,60,66, 68,71	B		High	RB(12.0 µg/L)
		57,60,66,68,71	B		High	RB(12.0 µg/L) CRH(120%,120%)
	MCRS73	73,74,77-79,82,83	B		High	RB(12.0 µg/L)
		75,76,84-89,91	B		High	RB(12.0 µg/L) CRH(120%,120%)
Ca	MCRS57	60,66	B		High	RB(1300 µg/L)
		70,81	B		High	PB(344 µg/L)
	MCRS73	75,88-91	B		High	RB(1300 µg/L)
Cr	MCRS57	70,81		UL	Low	CRL(80.0%)
Cu	MCRS57	81	B		High	CCB(14.0 µg/L) CRH(140%,124%)
	MCRS73	88,90,91	B		High	RB(12.0 µg/L) CRL(78.0%,84.0%)
		89	B		High	RB(12.0 µg/L)
Pb	MCRS73	88-90	B		High	RB(51.2 µg/L)

+ All samples begin with the prefix MCRS.

* See explanation of Comments on Table 1B.

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TABLE 1A
SUMMARY OF QUALIFIERS ON DATA SUMMARY
FORM AFTER DATA VALIDATION

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ANALYTE	SDG	SAMPLES AFFECTED+	POSITIVE VALUES	NON- DETECTED VALUES	BIAS	COMMENTS*
Hg	MCRS57	57,59-66,69,71	B		High	RB(0.5 µg/L) MSH(130%)
		70,81	B		High	PB(0.2 µg/L)
		58,67,68,72	K		High	MSH(130%)
	MCRS73	73,75-80,82,85,90, 91	B		High	RB(0.50 µg/L)
Ni	MCRS57	57,60,62,66,68,71	B		High	RB(19.0 µg/L)
		70,81	B		High	CCB(23.0 µg/L) CRH(118%,126%)
	MCRS73	80,84-91	B		High	RB(19.0 µg/L) CRH(126%,114%)
		76,78,82,83	B		High	RB(19.0 µg/L)
K	MCRS57	All soil samples	B		High	PB(642 mg/Kg)
		70,81		UL	Low	CCN(-4590 µg/L)
Se	MCRS57	61,64,66,67,69,72	K		High	MSH(159%) CRH(136%)
	MCRS73	All samples except 79,80,86		R	Extr. Low	CRE(18.0%)
		79,80,86	J			MIX(127%,18.0%)
Na	MCRS57	All soil samples	B		HIGH	RB(1070 µg/L)
		70,81	B		HIGH	PB(718 µg/L)
	MCRS73	All samples	B		HIGH	PB(286 mg/Kg)
Tl	MCRS57	70		UL	LOW	ANL(83.0%)
	MCRS73	73,76-78,80,82-84, 88		UL	LOW	CCN(-6.0 µg/L)
Zn	MCRS57	57,60-63,65-69,71, 72	B		High	RB(2140 µg/L)
	MCRS73	All samples	B		High	RB(2140 µg/L)

+ All samples begin with the prefix MCRS.
* See explanation of Comments on Table 1B.

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TABLE 1B

CODES USED IN COMMENTS COLUMN

- CCN - The continuing calibration blank had a negative result with absolute value greater than the IDL (the result is in parenthesis). The quantitation limits may be biased low.
- CRL - The CRDL standard recovery was low (< 90%) [% recovery is in parenthesis]. The reported results which are < 2x CRDL and quantitation limits may be biased low.
- MSL - The matrix spike recovery was low (30% - 75%) [% recovery is in parenthesis]. The reported results and quantitation limits may be biased low.
- CRE - The CRDL standard recovery was extremely low (< 50%) [% recovery is in parenthesis]. The quantitation limits are unusable.
- RB - The rinsate blank had a result > IDL (the result is in parenthesis). The reported results which are < 5x the blank concentration may be biased high.
- CRH - The CRDL standard recovery was high (> 110%) [% recovery is in parenthesis]. The reported results which are < 2x CRDL may be biased high.
- CCB - The continuing calibration blank had a result > IDL (the result is in parenthesis). The reported results which are < 5x the blank concentration may be biased high.
- PB - The preparation blank had a result > IDL (the result is in parenthesis). The reported results which are < 5x the blank concentration may be biased high.
- MIX - The matrix spike recovery was high (> 125%) and the CRDL standard recovery was extremely low (< 50%) [% recoveries are in parenthesis]. The reported results which are < 2x CRDL estimated.
- ANL - The analytical spike recovery was low (< 85%) [% recovery is in parenthesis]. The quantitation limits may be biased low.

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TABLE 2

GLOSSARY OF DATA QUALIFIER CODES (INORGANIC)

CODES RELATED TO IDENTIFICATION

(confidence concerning presence or absence of analytes):

U = Not detected. The associated number indicates approximate sample concentration necessary to be detected.

(NO CODE) = Confirmed identification.

B = Not detected substantially above the level reported in laboratory or field blanks.

R = Unusable result. Analyte may or may not be present in the sample. Supporting data necessary to confirm result.

CODES RELATED TO QUANTITATION

(can be used for both positive results and sample quantitation limits):

J = Analyte Present. Reported value may not be accurate or precise.

K = Analyte present. Reported value may be biased high. Actual value is expected to be lower.

L = Analyte present. Reported value may be biased low. Actual value is expected to be higher.

[] = Analyte present. As values approach the IDL the quantitation may not be accurate.

UJ = Not detected, quantitation limit may be inaccurate or imprecise.

UL = Not detected, quantitation limit is probably higher.

OTHER CODES

Q = No analytical result.

+ = Result reported from dilution. Quantitation limit is elevated.

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TABLE 4

SAMPLES EXCEEDING THE CHEMICAL HEALTH ADVISORY LEVELS

SDG MCRS57
($\mu\text{g/L}$)

<u>Analyte</u>	<u>Advisory Limit</u>	<u>Sample</u>	<u>Concentration</u>
Cd	8	MCRS70	12.0
Pb	20	MCRS70	35.1
		MCRS81	51.2

SDG MCRS57
(mg/Kg)

<u>Analyte</u>	<u>Advisory Limit</u>	<u>Sample</u>	<u>Concentration</u>
Pb	500	MCRS64	595
		MCRS65	785
		MCRS67	698
		MCRS69	698
		MCRS72	1120

SDG MCRS73
(mg/Kg)

<u>Analyte</u>	<u>Advisory Limit</u>	<u>Sample</u>	<u>Concentration</u>
Pb	500	MCRS73	525
		MCRS74	567
		MCRS77	3380
		MCRS78	2310
		MCRS79	870
		MCRS82	530
		MCRS83	817

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TABLE 3

FIELD DUPLICATE RESULTS OUTSIDE CRITERIA

Criteria for soil samples: 35% RPD, \pm 2x CRDLSDG: MCRS73

<u>Analyte</u>	<u>Concentration (mg/Kg)</u>		<u>RPD</u>
	<u>MCRS77</u>	<u>MCRS78</u>	
Ba	894	542	49.0
Ca	14200	31300	75.2
Mg	3890	7850	3960+

RPD = Relative Percent Difference

+ = Difference in mg/Kg instead of RPD.

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WATER SAMPLES

Case #: 22533 Sampling Date(s): 08/03/94

♦ Due to dilution, sample quantitation limit is affected
See dilution table for specifics

Sample No.	MCS870	MCS881
Dilution Factor	1.0	1.0
Location	B9-1	B9-2 RINSATE
CODL ANALYTE	RINSATE BLANK	RINSATE BLANK
200 Aluminum	UL	UL
60 Antimony	UL	UL
10 Arsenic	UL	UL
200 Barium		
5 Beryllium		
5 Cadmium	12.0	
5000 Calcium	(531)	B (1300)
10 Chromium	UL	UL
50 Cobalt		
25 Copper		(12.0) B
100 Iron	(77.0)	
3 Lead	35.1	51.2
5000 Magnesium		
15 Manganese	(10.0)	
0.2 Mercury	0.50 B	0.50 B
40 Nickel	(19.0) B	(15.0) B
5000 Potassium	UL	UL
5 Selenium		
10 Silver		
5000 Sodium	(874) B	(1070) B
10 Thallium	UL	
50 Vanadium		
20 Zinc	1560	2140
10 Cyanide		

CAOL = Contract Required Detection Limit

Action Level Exists

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

R310638

DATA SUMMARY FORM: INORGANICS
TABLE - 3
SOIL SAMPLES
(mg/kg)

Site Name: WESTINGHOUSE SHARON

Case #: 22533 Sampling Date(s): 08/02/94

SDG #: MCRS57

↑ Due to dilution, sample quantitation limit is affected
See dilution table for specifics

Sample No.	MCRS57	MCRS58	MCRS59	MCRS60	MCRS61	MCRS62	MCRS63	MCRS64	MCRS65	MCRS66
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
X Solids	93.0	79.5	91.6	78.2	93.6	79.3	80.3	78.2	87.5	81.8
Location	301	302	303	304	305	306	306 DUP	307	308	309
CDL ANALYTE						DUPLICATE OF	DUPLICATE OF			
						MCRS63	MCRS62			
40 Aluminum	13400	14400	17500	12300	13500	13300	13700	11800	10400	8850
12 Antimony	UL	UL	UL	UL	UL	UL	UL	UL	UL	UL
2 Arsenic	10.2	13.6	12.6	19.1	26.9	19.9	15.3	24.2	16.7	26.6
40 Barium	116	181	244	91.4	146	163	163	166	255	93.6
1 Beryllium	1.1	1.7	1.9		1.1	1.3	1.2	11.01	10.88	
1 Cadmium	1.9	3.9	3.9	2.0	3.6	4.3	3.9	4.8	2.4	2.0
1000 Calcium	12800	29800	36400	1470	6940	17000	17000	3170	3540	1360
2 Chromium	23.6	26.9	33.6	22.8	29.0	22.2	22.7	24.6	30.8	27.1
10 Cobalt	19.31	17.91	18.61	18.91	10.8	16.61	16.41	17.11	19.51	18.01
5 Copper	31.2	49.3	48.4	46.2	82.7	40.4	48.3	34.9	29.5	34.0
20 Iron	26400	27100	24800	27100	25800	23400	23300	24700	29000	28700
0.6 Lead	114	296	383	422	347	260	249	595	785	440
1000 Magnesium	3290	3350	6050	2150	2500	3250	3210	1890	2260	1960
3 Manganese	867	1560	1210	755	1130	958	1020	618	681	616
0.1 Mercury	0.16	1.1	0.33	0.23	0.21	0.13	0.31	0.38	0.23	0.43
8 Nickel	20.0	24.7	24.8	20.4	23.7	23.7	24.7	26.6	23.3	18.9
1000 Potassium	1940	1960	2200	2030	1630	1800	1770	110501	1260	19251
1 Selenium					1.4	K		1.5	K	1.3
2 Silver										
1000 Sodium	12663	15251	14001	11951	12311	12951	13131	12061	12591	12271
2 Thallium										
10 Vanadium	29.9	21.0	22.7	31.3	26.4	19.7	23.4	23.5	25.9	25.3
4 Zinc	913	3400	2530	854	1750	1940	1850	2840	1500	1190
0.5 Cyanide	0.45	0.63	0.71	0.83	1.0	0.63		0.64	1.0	0.73

CDL = Contract Required Detection Limit

*Action Level Exceeds

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR300639

DATA SUMMARY FORM: INORGANICS

Page 3 of 5

TABLE - 3

SOIL SAMPLES

(mg/Kg)

Site Name: WESTINGHOUSE SAISON

Case #: 22533 Sampling Date(s): 08/02/94 - 08/03/94

SDG #: NCSS57

+ Due to dilution, sample quantitation limit is affected
See dilution table for specifics

Sample No.	NCSS67	NCSS68	NCSS69	NCSS71	NCSS72
Dilution Factor	1.0	1.0	1.0	1.0	1.0
% Solids	88.8	94.1	76.9	81.2	87.2
Location	310	311	312	313	314
CADL ANALYTE					
40 Aluminum	9790	11000	14400	12400	17800
12 Antimony		UL	UL	UL	UL
2 Arsenic	24.9	16.6	34.9	16.9	19.4
40 Barium	223	259	248	86.3	328
1 Beryllium	10.391	10.853	2.3		2.1
1 Cadmium	3.8	2.1	4.9	1.7	3.9
1000 Calcium	6150	3530	9300	3170	34200
2 Chromium	39.9	21.3	60.8	22.2	30.3
10 Cobalt	17.61	17.91	16.2	19.01	18.31
5 Copper	109	64.4	92.2	21.7	73.2
20 Iron	35000	26600	36100	24300	27100
0.6 Lead	698	476	698	156	1120
1000 Magnesium	1960	2130	1510	2270	5230
3 Manganese	760	635	670	673	1600
0.1 Mercury	0.79	0.58	0.26	0.25	0.76
8 Nickel	25.4	19.1	34.5	21.5	23.4
1000 Potassium	1230	1330	1330	1920	1790
1 Selenium	2.1	K	2.0	K	1.4
2 Silver					
1000 Sodium	13061	12871	14891	12951	15571
2 Thallium					
10 Vanadium	23.6	23.6	42.5	23.9	23.9
4 Zinc	1800	1080	2180	897	1360
0.5 Cyanide	1.1		1.0		0.75

CADL = Contract Required Detection Limit

*Action Level Exceeds

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR300640

DATA SUMMARY FORM: INORGANICS

TABLE - 3

SOIL SAMPLES

(mg/kg)

Site Name: WESTINGHOUSE SHARON

Case #: 22533 Sampling Date(s): 08/03/94

SOG #: MCRS73

→ Due to dilution, sample quantitation limit is affected
See dilution table for specifics.

Sample No.	MCRS73	MCRS74	MCRS75	MCRS76	MCRS77	MCRS78	MCRS79	MCRS80	MCRS82	MCRS83
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
% Solids	88.4	88.3	89.0	79.8	85.3	85.7	71.9	83.1	83.7	85.3
Location	S15	S16	S17	S18	S19	S19 DUP	S20	S21	S22	S23
					DUPLICATE OF					
CDL ANALYTE					MCRS78	MCRS77				
40 Aluminum	12000	11500	11500	8830	8670	10400	11500	10300	13800	26400
12 Antimony										
2 Arsenic	26.9	28.6	23.4	18.8	21.9	18.3	18.8	29.0	21.2	19.8
40 Barium	156	157		143	994	542	375	194	700	646
1 Beryllium	1.3	1.1		10.991	10.881	1.1	11.1	1.7		
1 Cadmium	3.4	3.6	1.1	2.2	4.2	4.6	4.4		6.5	6.1
1000 Calcium	2560	2710	112301	4910	14200	31300	10500	31000	7960	14000
2 Chromium	26.0	27.1	18.5	26.1	43.6	34.3	40.2	19.3	31.1	33.0
10 Cobalt	19.31	19.01	19.81	18.41	18.81	17.31	18.81	13.91	11.01	110.01
5 Copper	38.9	40.1		48.1	42.5	40.3	71.3	18.5	93.9	100
20 Iron	30700	27100	36000	22300	29200	23300	28100	13600	23900	22700
0.6 Lead	525	567	105	353	3380	2310	870	96.3	530	817
1000 Magnesium	1950	1820	123601	1850	3890	7850	2550	5440	2560	5300
3 Manganese	816	841	517	521	1090	881	846	570	843	938
0.1 Mercury	0.51	1.1	0.51	0.54	0.21	0.29	0.35	0.12	0.36	0.76
8 Nickel	21.8	22.0	22.3	22.1	23.0	20.1	36.9	15.4	22.0	19.8
1000 Potassium	1140	18281	1430	19731	18781	19521	112801			
1 Selenium										
2 Silver										
1000 Sodium	13661	13771	13451	14221	14291	15101	16471	16261	14801	14421
2 Thallium										
10 Vanadium	27.3	25.8	26.0	19.9	17.9	17.2	27.5	14.0	23.7	24.6
4 Zinc	1970	2160	206	1370	1270	1240	2230	205	1700	1690
0.5 Cyanide					0.59	0.76			0.78	
CDL = Contract Required Detection Limit										
Action Level Exceeds										
SEE NARRATIVE FOR CODE DEFINITIONS										

revised 07/90

AR300641

DATA SUMMARY FORM: INORGANICS

TABLE - 3

SOIL SAMPLES

(mg/Kg)

Page 5 of 5

Site Name: WESTINGHOUSE SHAWON

Case #: 22533 Sampling Date(s): 04/03/94

SDG #: MCR575

+ Due to dilution, sample quantitation limit is affected.
See dilution table for specifics.

Sample No.	MCR584	MCR585	MCR586	MCR587	MCR588	MCR589	MCR590	MCR591
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
% Solids	90.0	84.6	84.7	86.0	79.6	87.0	80.2	75.9
Location	S24	S25	S26	S27	S28	S29	S30	S31
CRCL ANALYTE								
40 Aluminum	7900	9490	6240	13200	10500	11700	12200	15700
12 Antimony								
2 Arsenic	9.6	13.0	23.2	11.2	18.3	12.9	10.4	12.2
40 Barium	63.5	101	97.3	122	147.21	68.0	56.8	60.1
1 Beryllium			1.2	1.2				
1 Cadmium	1.2	1.8	10.94	1.2	11.0	10.92		11.1
1000 Calcium	4840	4180	3280	19300	1460	1963	1491	311
2 Chromium	20.2	22.7	14.4	16.7	20.6	15.2	14.4	16.6
10 Cobalt	15.8	17.1	17.8	17.2	110.3	19.4	17.1	13.7
5 Copper	18.7	34.9	31.4	27.0	10.8	13.1	10.5	9.0
20 Iron	25200	30800	28600	22100	21500	22200	19200	19500
0.6 Lead	360	190	202	223	40.4	51.8	44.4	74.5
1000 Magnesium	1750	1870	1850	3700	3390	2080	1730	1730
3 Manganese	603	581	404	1120	676	710	369	1340
0.1 Mercury		0.24	1.7				0.12	0.13
8 Nickel	16.3	18.6	17.5	17.7	19.1	18.4	14.4	17.7
1000 Potassium	1730	1700			1430	1340	1250	1110
1 Selenium			1.7					
2 Silver								
1000 Sodium	1317	1349	1449	1674	1420	1333	1445	1451
2 Thallium								
10 Vanadium	16.3	23.9	26.4	20.7	20.6	22.5	22.5	27.1
4 Zinc	758	877	212	809	120	147	166	136
0.5 Cyanide			0.94	0.93		0.80		

CRCL = Contract Required Detection Limit

*Action Level Exceeds

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR300642



United States Environmental Protection Agency
Contract Laboratory Program Sample Management Office
PO Box 918 Alexandria, VA 22313
703-557-2490 FTS 557-2490

Inorganic Traffic Report & Chain of Custody Record

(For Inorganic CLP Analysis)

SAS No.
(If applicable)

Case No.
22553

1. Project Code 5303		2. Region No. III		3. Sampling Co. WESTON		4. Date Shipped 8/4/94		Carrier Federal Express		5. Ship To 202 868 0964		6. Preservative (Enter in Column D) 1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2Cr2O7 6. Ice only 7. Other N. Not preserved		7. Sample Description (Enter in Column A) 1. Surface Water 2. Ground Water 3. Leachate 4. Rinsate 5. Soil/Sediment 6. Oil (High only) 7. Waste (High only) 8. Other (Specify)		8. Enter Appropriate Qualifier for Designated Field QC B - Blank S - Spike D - Duplicate PE - Perform Eval. -- = Not a QC Sample			
Regional Information		Sampler (Name) J. E. CARTER		Airbill Number 202 868 0964		ATTN: John Mack, Mgr		Mack Laboratories, Inc 2199 Dartmore Avenue Pittsburgh, PA 15210		Sample Initials JL		Corresp. CLP Org. Samp. No. COR13		Mo/Day/Year/Time Sample Collection 08/02/94 11:15		Station Location Number S01		Regional Specific Tracking Number or Tag Numbers T3-22035	
Non-Superfund Program		Sampler Signature J. E. Carter		Removal SF <input checked="" type="checkbox"/> RIFS <input type="checkbox"/> CLEM PRP <input type="checkbox"/> REMA ST <input type="checkbox"/> REM FED <input type="checkbox"/> OIL LSI <input type="checkbox"/> UST		3. Type of Activity PA <input type="checkbox"/> RA <input type="checkbox"/> O&M SS <input type="checkbox"/> NPLD		E - RAS Analysis Metallic Total Dissolved Cyanide Nitrite Nitrate Fluoride pH Conductivity		D Preservative vial from Box 6		C Sample Type: Med Comp./ Grab		B Conc. Low Med High		A Enter from Box 7		CLP Sample Numbers (from labels)	
Site Name WESTINGHOUSE SHARON		City, State Sharon, PA		Site Spill ID		CLP Sample Numbers (from labels)		MCRS57		5		Low		Grab		5		MCRS57	
								MCRS58		5		Low		Grab		5		MCRS58	
								MCRS59		5		Low		Grab		5		MCRS59	
								MCRS60		5		Low		Grab		5		MCRS60	
								MCRS61		5		Low		Grab		5		MCRS61	
								MCRS62		5		Low		Grab		5		MCRS62	
								MCRS63		5		Low		Grab		5		MCRS63	
								MCRS64		5		Low		Grab		5		MCRS64	
								MCRS65		5		Low		Grab		5		MCRS65	
								MCRS66		5		Low		Grab		5		MCRS66	
Shipment for Case complete? (Y/N)		Page 1 of 4		Sample used for a spike and/or duplicate		MCRS62		Additional Sampler Signatures J. Mack		Chain of Custody Seal Number									

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature) J. E. Carter	Date / Time 8/4/94 1600	Received by: (Signature)	Date / Time	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
Remarks		Is custody seal intact? Y/N/none				

EPA Form 9110-1 (Rev. 5-91) Replaces EPA Form (2075-6), previous edition which may be used
DISTRIBUTION:
Green - Region Copy Pink - SMO Copy White - Lab Copy for return to Region Yellow - Lab Copy for return to SMO

Split Samples ☐ Accepted ☐ Declined

AR300643

Inorganic Traffic Report & Chain of Custody Record

SAS No.
(if applicable)

Case No.

[illegible]

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time	Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time	Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time
<i>Chen V. A. L.</i>	5/14/94 1600										
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time	Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time	Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time
Relinquished by: (Signature)	Date / Time	Received for Laboratory by: (Signature)	Date / Time	Remarks	Is custody seal intact?	Y/N/None					

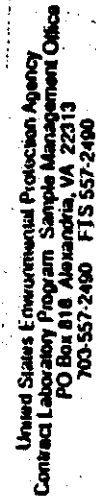
EPA Form 910-1 (Rev. 5-01) Replaces EPA Form (2075-0), previous edition which may be used

DISCLAIMER

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Copy for Re. SMO
/ Pink - SMO Copy
White - Lab. Copy for return to Region
Yellow - Lab

SEE REVERSE FOR ADDITIONAL STANDARD INSTRUCTIONS

AR300644



Inorganic Traffic Report & Chemical Analysis of Custody Record

Case No.

SAS No.
(if applicable)

Inorganic Traffic Report & Check of Custody Record (For Inorganic CLP Analysis)

United States Environmental Protection Agency
Contract Laboratory Program Sample Management Office
PO Box 818 Alexandria, VA 22313
703-557-2490 FIS 557-2490

1. Project Code	Account Code	2. Region No.	Sampling Co.	4. Date Shipped	Carrier	7. Sample Description (Enter in Column A)
5303		III	WESTON	8/14/94	Federal Express	1. Surface Water 2. Ground Water 3. Leachate 4. Rinsate 5. Soil/Sediment 6. Oil (High only) 7. Waste (High only) 8. Other (Specify)
Regional Information		3. Type of Activity	5. Ship To	6. Preservative (Enter in Column D)	8. Other (Specify)	9. Enter Appropriate Qualifier for Designated Field QC
Non-Superfund Program		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>	Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.	1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved		K Enter Appropriate Qualifier for Designated Field QC B = Blank S = Spike D = Duplicate PE = Perform. Eval. -- = Not a QC Sample
Site Name		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
WESTINGHOUSE - SHARON		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
City, State		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
SHARON, PA		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
Site Name		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
WESTINGHOUSE - SHARON		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
City, State		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
SHARON, PA		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
Site Name		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
WESTINGHOUSE - SHARON		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
City, State		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
SHARON, PA		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
Site Name		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
WESTINGHOUSE - SHARON		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
City, State		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
SHARON, PA		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
Site Name		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
WESTINGHOUSE - SHARON		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
City, State		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
SHARON, PA		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
Site Name		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
WESTINGHOUSE - SHARON		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
City, State		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
SHARON, PA		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <input checked="" type="checkbox"/> SS <input checked="" type="checkbox"/> LS <input checked="" type="checkbox"/> OSM <input checked="" type="checkbox"/> NPLD <input checked="" type="checkbox"/> OIL <input checked="" type="checkbox"/> UST <input checked="" type="checkbox"/>		Mack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr.		1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2CR2O7 6. Ice only 7. Other (Specify) N. Not preserved
Site Name		3. Type of Activity		5. Ship To		6. Preservative (Enter in Column D)
WESTINGHOUSE - SHARON		Lead <input checked="" type="checkbox"/> PFAS <input checked="" type="checkbox"/> PA <				

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature) <i>[Signature]</i>	Date / Time 8/4/94/1608	Received by: (Signature)	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Relinquished by: (Signature)	Date / Time	Received by: (Signature)
Relinquished by: (Signature)	Date / Time	Received for Laboratory by: (Signature)	Date / Time	Remarks	is custody seal intact? Y/N/none
EPA Form 9116-1 (Rev. 5-93) Replaces EPA Form (2075-9), previous edition which may be used			Split Samples <input checked="" type="checkbox"/> Accepted (Signature) <input type="checkbox"/> Declined		

DISTRIBUTION:
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AR300645

DISTRIBUTION:
Green - Region Copy Pink - SMO Copy White - Lab Copy for return to Region Yellow - Lab Copy for Return to SMO



United States Environmental Protection Agency
Contract Laboratory Program Sample Management Office
PO Box 618 Alexandria, VA 22313
703-557-2400 FTS 557-2400

Inorganic Traffic Report & Chain of Custody Record

(For Inorganic CLP Analysis)

SAS No.
(if applicable)

Case No.

22533

1. Project Code 5303	Account Code	2. Region No. III	Sampling Co. WESTON	4. Date Shipped 8/14/94	Carrier Federal Express	6. Preservative (Enter in Column D) 1. HCl 2. HNO3 3. NaOH 4. H2SO4 5. K2Cr2O7 6. Ice only 7. Other N. Not preserved	7. Sample Description (Enter in Column A) 1. Surface Water 2. Ground Water 3. Leachate 4. Runoff 5. Soil/Sediment 6. Oil (High only) 7. Waste (High only) 8. Other (Specify)				
Regional Information		Sample (Name) JOE CARTER		Airbill Number 2028680964							
Non-Superfund Program		Sample Signature <i>Joe S. Carter</i>		5. Ship To Hack Laboratories, Inc. 2199 Dartmore Avenue Pittsburgh, PA 15210 ATTN: John Mack, Mgr							
Site Name Westinghouse - Sharon		3. Type of Activity SF <input checked="" type="checkbox"/> PAFS <input type="checkbox"/> CLEM <input type="checkbox"/> PRP <input type="checkbox"/> PA <input type="checkbox"/> RD <input type="checkbox"/> REM <input type="checkbox"/> ST <input type="checkbox"/> SS <input type="checkbox"/> OAM <input type="checkbox"/> OIL <input type="checkbox"/> FED <input type="checkbox"/> LSI <input type="checkbox"/> NPL <input type="checkbox"/> UST <input type="checkbox"/>									
CLP Sample Numbers (from labels)	A Enter from Box 7	B Conc. Low Med High	C Sample Type: Comp/Grab	D Preservative from Box 6	E - RAS Analysis Metals: Described, Graded, Residue, Spill, Other, pH, TOC, etc.	F Regional Specific Tracking Number or Tag Numbers	G Station Location Number	H Mo/Day/Year Sample Collection	I Sampler Initials	J Corresp. CLP Org. Sample No.	K Enter Appropriate Qualifier for Designated Field QC B = Blank S = Spike D = Duplicate PE = Perform Eval. -- = Not a QC Sample
MCRS65	5	Low	Grab	6		T3-22002	525	08/03/94 1600	JC	MCR41	
MCRS66	5	Low	Grab	6		T3-22003	526	08/03/94 1600	JC	MCR42	
MCRS67	5	Low	Grab	6		T3-22004	527	08/03/94 1645	JC	MCR43	
MCRS68	5	Low	Grab	6		T3-22005	528	08/03/94 1735	JC	MCR44	
MCRS69	5	Low	Grab	6		T3-22006	529	08/03/94 1800	JC	MCR45	
MCRS90	5	Low	Grab	6		T3-22007	530	08/03/94 1845	JC	MCR46	
MCRS91	5	Low	Grab	6		T3-22008	531	08/03/94 1915	JC	MCR47	
Shipment for Case complete? (Y/N)	Page 4 of 4	Sample used for a spike and/or duplicate		Additional Sample Signatures <i>John Mack</i>							

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature) <i>Ch. V. Felt</i>	Date / Time 8/14/94 1600	Received by: (Signature)	Date / Time
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time
Remarks: Is custody seal intact? Y/N none			

EPA Form 910-1 (Rev. 5-91) Replaces EPA Form (2075-6), previous edition which may be used
DISTRIBUTION: Green - Re. Copy for Retention; Pink - SMO Copy; White - Lab Copy for return to Region; Yellow - Lab Copy for Retention to SMO

AR300646C

Attachment D

Organics Data Validation, Analytical Summary, Traffic Report

AR300647



Environmental Systems & Technologies Co.

Environmental Services Assistance Teams
Region 3
1419 Forest Drive, Suite 104
Annapolis, Maryland 21403

Phone: (410) 268-7705
Fax: (410) 268-8472

DATE: September 23, 1994

SUBJECT: Organic Data Validation for Case 22533
Site: Westinghouse Sharon

FROM: Edgar A. Latham ^{EL}
Senior Organic Data Reviewer
Mahmoud Hamid ^{MH}
Senior Oversight Chemist
Mahboobeh Mecanic ^{MM}
Senior Oversight Chemist

TO: Susanne Stevens
ESAT Acting Regional Project Officer

THROUGH: Dale S. Boshart ^{DB}
ESAT Team Manager

OVERVIEW

Case 22533 was submitted to EnviroSystems Laboratory (ENVSYS) for full organic analysis. The case consisted of thirty-three (33) soil samples, one (1) trip blank that was analyzed for volatiles only, and two (2) rinsate blanks. The samples were submitted in two (2) sample delivery groups (SDGs). One (1) soil field duplicate pair for each SDG was analyzed. The samples were analyzed as a Contract Laboratory Program (CLP) Routine Analytical Service (RAS).

SUMMARY

All samples were successfully analyzed for all target compounds except for 2,4-dinitrophenol. All other instrument and method sensitivities were according to the Contract Laboratory Program (CLP) Routine Analytical Service (RAS) protocol.

MAJOR PROBLEM

- o The average relative response factor for 2,4-dinitrophenol was less than 0.05 (<0.05) in the semivolatile continuing calibrations dated 8/15/94 and 9/1/94. The quantitation limit for this compound was qualified "R" for the associated samples. See Table I in Appendix E.

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- o The semivolatile reextraction of sample CQR19 was performed twenty (20) days after the date of sample collection. Although no technical holding time for the extraction of semivolatiles in soil samples has been established, the technical holding time of seven (7) days for aqueous samples has been exceeded by thirteen (13) days. The aqueous sample holding time was applied; therefore, the quantitation limits were qualified "UJ" and positive results were qualified "J", unless superseded by the "B" qualifier. The contractual holding time of ten (10) days from the validated time of sample receipt was exceeded by seven (7) days because the sample was reanalyzed due to surrogate outliers. See Form Is in Appendix C, Traffic Report, and Case Narrative in Appendix E.
- o Several compounds failed precision criteria in the volatile and semivolatile initial and/or continuing calibrations. The positive results were qualified "J", except when superseded by the "B" qualifier, and when the percent relative standard deviation (%RSD) or the percent difference (%D) was greater than fifty percent (>50%), the quantitation limits were qualified "UJ", except when superseded by the "R" qualifier for these compounds in the affected samples.
- o Benzo(k)fluoranthene exceeded the linear calibration range in the semivolatile analysis of sample CQR25. However, this compound was not detected in the diluted analysis. The result was reported from the initial analysis and qualified "J".
- o In the semivolatile analyses of samples CQR16RE, CQR19, CQR40RE, and CQR42 several compounds were detected in either the initial or reanalyses. These compound were marked with a plus sign (+) on the data summary forms (DSFs) and qualified "J", unless superseded by the "B" qualifier.
- o In the volatile analyses, samples CQR25, CQR32, and CQR42 had the system monitoring compound toluene-d8 outside the upper QC limits. Reanalyses of samples CQR25 and CQR42 produced similar results. The MS/MSD analyses of sample CQR32 confirmed matrix interference through internal standard responses. Quantitation limits were qualified "UJ". No positive results were detected.
- o The semivolatile base/neutral surrogates 2-fluorobiphenyl (S2) and terphenyl-d14 (S3), in addition to the acid surrogate 2,4,6-tribromophenol (S6) were outside the upper QC limits for sample CQR25. The positive results for the base/neutral compounds were qualified "K", unless superseded by the "J" qualifier. See Form II SV-2 in Appendix E.

- o In the pesticide/PCB analyses, samples CQR14, CQR23, CQR29, and CQR30 had two (2) surrogate recoveries outside the upper QC limits on different columns. Quantitation limits and reported results were qualified "UJ" and "J", respectively on the DSFs.
- o In the pesticide/PCB analyses, sample CQR21 had one (1) surrogate recovery outside the upper QC limits on one column and another surrogate recovery outside the lower QC limits on the other column. Quantitation limits and Positive results were qualified "UJ" and "J", respectively on the DSF.
- o In the pesticide/PCB analyses, samples CQR26, CQR39, CQR41, CQR44, CQR45, and CQR47 had two (2) surrogate recoveries outside the lower QC limits on different columns. Quantitation limits and reported results were qualified "UJ" and "J", respectively on the DSFs.
- o In the pesticide/PCB analyses, samples CQR42 and CQR46 had four (4) surrogate recoveries outside the lower QC limits on both columns. Positive results were qualified "L" and quantitation limits were qualified "UL" on the DSFs.
- o In the volatile analyses, samples CQR24, CQR29, CQR30, CQR31, CQR34, CQR35, CQR42, and CQR47 had one (1) or more internal standards outside the QC limits. Reanalyses of the above samples produced similar results for CQR29, CQR30, CQR31, CQR34 and CQR42 and better QC results for samples CQR24RE, CQR35RE and CQR47RE. Results from initial analyses of samples CQR29, CQR30, CQR31, CQR34 and CQR42, and from the reanalyses of samples CQR24, CQR35, and CQR47 were reported on the DSFs. Quantitation limits for compounds quantitated using any of the failed internal standards were qualified "UJ". No positive results were detected.
- o In the semivolatile analyses, samples CQR19, CQR42, and CQR47 had the internal standards chrysene-d12 (IS5) and/or perylene-d12 (IS6) outside the QC limits. The reanalysis of these samples revealed similar results. Results were reported from the initial analyses. The quantitation limits were qualified "UJ" and positive results were qualified "J" for the compounds quantitated using the affected internal standards. See Form VIII SV-2 in Appendix E.
- o In the semivolatile analyses, samples CQR13, CQR18, and CQR23 had internal standards phenanthrene-d10 (IS4) and/or IS5 outside the QC limits, while samples CQR25, CQR29, and CQR30 had IS6 outside the QC limits. These samples were diluted and reanalyzed because several compounds exceeded the linear calibration range. The internal standard recoveries for the diluted analyses were within QC limits. Quantitation limits were qualified "UJ" and positive results were qualified "J",

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unless superseded by the "B" qualifier for compounds reported from the initial analyses that are associated with the affected internal standards. See Form VIII SV-2 in Appendix E.

- o In the semivolatile analyses, samples CQR40 and CQR41 had internal standards IS5 and IS6 outside the QC limits. The reanalysis of these samples revealed only IS6 as an outlier. Results were reported from the reanalyses. The quantitation limits were qualified "UJ" and positive results were qualified "J" for the compounds quantitated using IS6. See Form VIII SV-2 in Appendix E.
- o Three (3) internal standards were outside the QC limits in the semivolatile analysis of sample CQR16. In the reanalysis of this sample, only internal standard IS5 was outside the QC limit. Results were reported from the reanalysis. The quantitation limits were qualified "UJ" and positive results were qualified "J" for the compounds quantitated using IS5. See Form VIII SV-2 in Appendix E.
- o The internal standard IS6 was outside the QC limit for the semivolatile analysis of sample CQR35DL. Positive results reported from the dilution were qualified "J" for the compounds quantitated using IS6. See Form VIII SV-2 in Appendix E.
- o The "P" qualifier used on the pesticide/PCB Form Is denotes a percent difference (%D) greater than 25 percent (>25%) between the reported results on the two (2) columns used for the analyses. These results were qualified "J" on the DSFs. See Form Is in Appendix C and Form Xs in Appendix E.

NOTES

- o The semivolatile extractions for samples CQR32-CQR36 and CQR38-CQR47 were performed eight (8) days after the date of sample collection. Although no technical holding time for the semivolatile extraction of soil samples has been established, the technical holding time of seven (7) days for aqueous samples has been exceeded by one (1) day. No data were qualified due to this minor holding time infraction because semivolatile compounds are considered persistent in soil samples. The contractual holding time was met by the laboratory. See Form Is in Appendix C and Traffic Report in Appendix E.
- o In the pesticide/PCB analyses, soil samples CQR13-CQR20, CQR22-CQR28, CQR30, and CQR35, had several compounds that were not confirmed by GC/MS. Compounds detected at a concentration greater than 330 µg/Kg must be confirmed by GC/MS according to the 3/90 Statement of Work (SOW).

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- o The maximum concentrations of all compounds found in the analyses of the trip, field, and laboratory method blanks are listed below. Samples with concentrations of common laboratory contaminants less than ten times (<10X) the blank concentration or with concentration of other contaminants less than five times (<5X) the blank concentration have been qualified "B" on the DSFs.

<u>Compound</u>	<u>Concentration</u>
Acetone *	14 µg/L
bis(2-ethylhexyl)phthalate *	3 J µg/Kg
Endosulfan I	0.016 J µg/L
4,4'-DDE	0.0064 J µg/L
Alpha-chlordane	0.018 J µg/L

* = Common laboratory contaminants

- o Two (2) field duplicate pairs (CQR18/CQR19) and (CQR33/CQR34) were analyzed. The results and precision estimates, excluding the blank contaminants, are listed below:

<u>Compound</u>	<u>Concentration (µg/Kg)</u>		<u>RPD</u>
	<u>CQR18</u>	<u>CQR19</u>	
2-methylnaphthalene	ND	79 J	IN
acenaphthylene	ND	82 J	IN
pentachlorophenol	ND	160 J	IN
phenanthrene	1400 J	1000 J	33
anthracene	300 J	200 J	40
carbazole	210 J	ND	IN
di-n-butylphthalate	180 J	820 J	128
fluoranthene	3600	1800 J	67
pyrene	2600	1500 J	54
benzo(a)anthracene	1800 J	1800 J	0
chrysene	1800 J	1200 J	40
benzo(b)fluoranthene	2900	1300 J	76
benzo(k)fluoranthene	2600	1300 J	67
benzo(a)pyrene	2000	1100 J	58
indeno(1,2,3-cd)pyrene	870	1700 J	65
dibenz(a,h)anthracene	510	1600 J	103
benzo(g,h,i)perylene	920 J	1400 J	41
4,4'-DDE	7.3 J	11 J	40
Endrin	18 J	20 J	10
4,4'-DDT	42 J	41 J	2
Endrin aldehyde	40 J	46 J	14
Alpha-Chlordane	51 J	60 J	16
Gamma-Chlordane	38 J	46 J	19
Aroclor-1260	1300	1400	7

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<u>Compound</u>	<u>Concentration ($\mu\text{g/Kg}$)</u>		<u>RPD</u>
	<u>COR33</u>	<u>COR34</u>	
acenaphthylene	82 J	95 J	15
phenanthrene	540	770	35
anthracene	230 J	250 J	8
carbazole	150 J	220 J	38
di-n-butylphthalate	120 J	ND	IN
fluoranthene	2100	2400	13
pyrene	2100 J	1800 J	15
benzo(a)anthracene	1400	1600	13
chrysene	1300	1500	14
benzo(b)fluoranthene	ND	2200	IN
benzo(k)fluoranthene	2300	ND	IN
benzo(a)pyrene	790	1200	41

<u>Compound</u>	<u>Concentration ($\mu\text{g/Kg}$)</u>		<u>RPD</u>
	<u>COR33</u>	<u>COR34</u>	
Endosulfan I	17 J	15 J	12
4,4'-DDE	9 J	12	28
4,4'-DDT	16 J	19 J	17
Endrin aldehyde	12 J	11 J	9
Alpha-Chlordane	14 J	17 J	19
Gamma-Chlordane	9.5	10	5
Aroclor-1260	240 J	220 J	9

RPD = Relative Percent Difference

ND = Not Detected

IN = Indeterminate

- o Many of the samples required dilutions for the semivolatile analyses because target compounds exceeded the linear calibration range. These compounds were qualified with an "E" by the laboratory. Results reported from the diluted analyses were marked with an asterisk (*) on the DSFs. The analyses of samples COR14 and COR28 were performed initially at a 10X dilution. See Form Is in Appendix C and the Case Narrative in Appendix E.
- o In the pesticide/PCB analyses, several compounds exceeded the linear calibration range in samples COR13 and COR16 and dilutions were performed. Results reported from the diluted analyses were marked with an asterisk (*) on the DSFs.
- o In the pesticide/PCB analysis of sample COR16, dieldrin was detected at a concentration that exceeded the linear calibration range in the initial analysis. This compound was not reported by the laboratory in the 20X dilution. However, an inspection of the raw data shows the compound was present. The result was calculated by the reviewer and the Form I corrected. Dieldrin

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was reported from the diluted analysis and marked with an asterisk (*) on the DSF. See Form Is in Appendix C and the chromatograms in Appendix E.

- o The laboratory indicated that samples were diluted during the pesticide/PCB analyses because high target compound levels. Sample CQR24 was analyzed at a 20X dilution, sample CQR27 at a 10X dilution, and samples CQR29 and CQR30 at a 2X dilution. The laboratory diluted and analyzed several other samples, however, in the reviewer's judgement, many of these dilutions were unnecessary because the raw data did not suggest the results exceeded the linear calibration range. See Form Is in Appendix C and Case Narrative in Appendix E.
- o Gel Permeation Chromatography (GPC) cleanup was performed in the semivolatile and pesticide/PCB analyses of the soil samples. The dilution factor of two (2) required by this procedure was accounted for in the analytical procedures used by the laboratory.
- o Non-spiked compounds, other than blank contaminants, were determined in samples CQR18, CQR32, and the MS/MSD analyses of these samples. The results and precision estimates are as follows:

<u>Compound</u>	<u>CQR18</u>	<u>CQR18MS</u>	<u>CQR18MSD</u>	<u>1RSD</u>
2-methylnaphthalene	ND	ND	130 J	IN
phenanthrene	1400 J	1000	1400	18
anthracene	300 J	240 J	300 J	12
carbazole	210 J	170 J	200 J	11
di-n-butylphthalate	180 J	180 J	200 J	6
fluoranthene	3600	3600 E	4500 E	13
benzo(a)anthracene	1800 J	1500	1800	10
chrysene	1800 J	1400	1800	14
benzo(b)fluoranthene	2900	2400	2600	10
benzo(k)fluoranthene	2600	1400	2400	30
benzo(a)pyrene	2000	1600	1800	11
indeno(1,2,3-cd)pyrene	870	850	790	5
dibenz(a,h)anthracene	510	ND	300 J	95 +
benzo(g,h,i)perylene	920 J	880	890	2
4,4'-DDE	7.3 J	ND	ND	IN
Endrin aldehyde	40 J	33 J	6.9 J	63
Alpha-chlordane	51 J	47 J	8.6 J	66
Gamma-chlordane	38 J	35 J	6.7 J	65
Aroclor-1260	1300	1100	240	64

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<u>Compound</u>	<u>CQR32</u>	<u>CQR32MS</u>	<u>CQR32MSD</u>	<u>%RSD</u>
phenanthrene	970	1000	1000	2
anthracene	160 J	190 J	200 J	11
carbazole	250 J	250 J	250 J	0
di-n-butylphthalate	790	850	920	8
fluoranthene	2700	2800	2800	2
benzo(a)anthracene	1300	990	1000	16
chrysene	980	1300	1200	14
benzo(b)fluoranthene	1700	ND	ND	IN
benzo(k)fluoranthene	ND	2600	2600	0 +
benzo(a)pyrene	200 J	660	780	56
indeno(1,2,3-cd)pyrene	420	ND	ND	IN
dibenz(a,h)anthracene	230 J	ND	ND	IN
benzo(g,h,i)perylene	160 J	ND	ND	IN
Endosulfan I	18	21	17 J	11
4,4'-DDE	6 J	7.2 J	6.5 J	9
Endrin aldehyde	14 J	14 J	13 J	4
Alpha-chlordane	3.6 J	3.8 J	3.3 J	7
Aroclor-1260	250 J	260 J	220 J	9

%RSD = Percent Relative Standard Deviation

+ = Relative Percent Difference (RPD) instead of %RSD

ND = Not Detected

IN = Indeterminate

E = Result exceeded the linear calibration range

- o In the MS/MSD volatile analyses of samples CQR18 and CQR32, one (1) out of ten (10) spike recoveries were outside the QC limits.
- o In the MS/MSD semivolatile analyses of sample CQR18, one (1) out of eleven (11) relative percent differences (RPDs) and two (2) out of twenty-two (22) spike recoveries were outside the QC limits.
- o In the MS/MSD semivolatile analyses of sample CQR32, two (2) out of eleven (11) RPDs and five (5) out of twenty-two (22) spike recoveries were outside the QC limits.
- o In the MS/MSD pesticide/PCB analyses of sample CQR18, six (6) out of six (6) RPDs and six (6) out of twelve (12) spike recoveries were outside the QC limits.
- o In the semivolatile analyses, several samples had one (1) or two (2) surrogate recoveries outside the upper QC limits. No data were qualified. Surrogate recoveries for diluted samples were designated with a "D" by the laboratory. The extraction efficiency/method accuracy cannot be verified. See Form II SV-1&2.

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- o In the pesticide/PCB analyses, several samples had one (1) surrogate recovery outside the QC limits. No action was taken.
- o In the pesticide/PCB analyses, the MSD analysis of sample CQR18 had four (4) surrogate recoveries outside the lower QC limits on both columns. No action was taken.
- o In the pesticide/PCB analyses, the surrogate decachlorobiphenyl (DCB) had a retention time outside the established window for several samples because of matrix interference. No action was taken.
- o The retention times for internal standards IS5 and IS6 were outside the upper QC limits in the semivolatila analysis of sample CQR35. Results for the associated compounds were reported from the dilution and marked with an asterisk (*).
- o Several internal standards were outside the QC limits in the semivolatila MS/MSD analyses of samples CQR18 and CQR32. No data were qualified.
- o The tentatively identified compounds (TICs) in Appendix D were reviewed and corrected during data validation. Compounds identified as blank contaminants, target compounds of another fraction, aldol condensation products, or phthalates were crossed off the TIC Form Is.

All data for Case 22533 were reviewed according to the National Functional Guidelines for Evaluating Organic Analyses with modification for use within Region III. The text of the report addresses only those problems affecting usability.

ATTACHMENTS

- 1) Appendix A - Glossary of Data Qualifiers
- 2) Appendix B - Data Summary Forms. These include:
 - (a) All positive results for target compounds with qualifier codes where applicable.
 - (b) All unusable detection limits (qualified "R").
- 3) Appendix C - Results as Reported by the Laboratory for All Target Compounds
- 4) Appendix D - Reviewed and Corrected Tentatively Identified Compounds
- 5) Appendix E - Support Documentation

DCN:EL409A01.WES

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GLOSSARY OF DATA QUALIFIER CODES (ORGANIC)

CODES RELATED TO IDENTIFICATION

(confidence concerning presence or absence of compounds)

- U = Not detected. The associated number indicates approximate sample concentration necessary to be detected.
- NO CODE = Confirmed identification.
- B = Not detected substantially above the level reported in laboratory or field blanks.
- R = Unusable result. Analyte may or may not be present in the sample. Supporting data necessary to confirm result.
- N = Tentative identification. Consider present. Special methods may be needed to confirm its presence or absence in future sampling efforts.

CODES RELATED TO QUANTITATION

(can be used for both positive results and sample quantitation limits):

- J = Analyte present. Reported value may not be accurate or precise.
- K = Analyte present. Reported value may be biased high. Actual value is expected to be lower.
- L = Analyte present. Reported value may be biased low. Actual value is expected to be higher.
- UJ = Not detected, quantitation limit may be inaccurate or imprecise.
- UL = Not detected, quantitation limit is probably higher.

OTHER CODES

- NJ = Qualitative identification questionable due to poor resolution. Presumptively present at approximate quantity.
- Q = No analytical result.

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STAINING WATER SAMPLES

(7/8)

**To calculate sample quantitation limits:
(EQL = Dilution Factor)**

Sample No.	Cor12	Cor26	Cor37
Dilution Factor	1.0	1.0	1.0
Location	TR-1	AB-1	AB-2
CBL	CONPOUND	SAMPLE IS A TRIP BLANK.	SAMPLE IS A RINSEATE BLANK.
10	Chloromethane		
10	Bromomethane		
10	Vinyl Chloride		
10	Chloroethane		
10	Methylene Chloride		
10	Acetone		
10	Carbon Dioxide		
10	1,1-Dichloroethane		
10	1,1-Dichloroethane		
10	Total 1,2-Dichloroethane		
10	Chloroform		
10	1,2-Dichloroethane		
10	n-Butane		
10	1,1,1-Trichloroethane		
10	Carbon Tetrachloride		
10	Bromodichloromethane		

CRQL = Contract Required Quantitation Limit**Action Level Exits**

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

AR300658

WATER SAMPLES

(7/6n)

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94

To calculate sample quantitation limits:
(CNDL • Dilution Factor)

[illegible]**CRCL = Contract Required Quantitation Limit**

***Action Level Exists**

SEE NARRATIVE FOR CODE DEFINITIONS

Revised 07/90

AK-100659

(wa/Ka)

To calculate sample quantitation limits:
 $(\text{CAL} \times \text{Dilution factor}) / ((100 - \text{Moisture})/100)$

[illegible]

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

AR:110660

SOIL SAMPLES

(b)(7)(D)

To calculate sample quantitation limits:
 $(CROL * Dilution factor / ((100 - \%moisture)/100))$

Sample No.	COR13	COR14	COR15	COR16	COR17	COR18	COR19	COR20	COR21
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
% Moisture	7	21	10	23	8	19	15	21	12
Location	301	302	303	304	305	306	306 DUPE	307	308
CRCL									
COMPOUND									
10	1,2-Dichloropropane								
10	Cis-1,3-Dichloropropene								
10	Trichloroethene								
10	Dibromochloroethane								
10	1,1,2-Trichloroethane								
10	Benzene								
10	Trans-1,3-Dichloropropene								
10	Bromoform								
10	4-Methyl-2-pentanone								
10	2-Hexanone								
10	Tetrachloroethene								
10	1,1,2,2-Tetrachloroethane								
10	Toluene								
10	Chlorobenzene								
10	Ethylbenzene								
10	Styrene								
10	Total Xylenes								

CRQL - Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

66/20 revised

AR300661

SOIL SAMPLES

(b)(7)(D)

To calculate sample quantitation limits:
 $(CROL * \text{dilution factor} / ((100 - Xmoisture)/100))$

[illegible]**CRQL = Contract Required Quantitation Limit**

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AK-100665

DATA SUMMARY FORM: VOLATILES 1

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Site Name: Westinghouse Sharon

SOIL SAMPLES
(ug/Kg)

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94

To calculate sample quantitation limits:
(CRL = Dilution factor / ((100 - Moisture)/100))

Sample No.	COR42	COR43	COR44	COR45	COR46	COR47NE
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0
% Moisture	13	16	20	12	20	25
Location	S26	S27	S28	S29	S30	S31
COMPOUND						
Chloroethane						
Bromoethane						
Vinyl Chloride						
Chloroethene						
Methylene Chloride						
Acetone						
Carbon Disulfide						
1,1-Dichloroethane						
1,1-Dichloroethane						
Total 1,2-Dichloroethane						
Chloroform						
1,2-Dichloroethane						
2-Butanone						
1,1,1-Trichloroethane						
Carbon Tetrachloride						
Bromodichloromethane						

CRL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

AR300666

Site Name: Westinghouse Sharon

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94

(b)(7)(C)
SOIL SAMPLES

To calculate sample quantitation limits:
 $(CROL * Dilution factor / ((100 - \%moisture)/100))$

[illegible]**CRQL = Contract Required Quantitation Limit**

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

44-110667

Site Name: Westinghouse Sharon

WATER SAMPLES
(1/20)

Case #: 22533 Sampling Date(s): 8-2-94 • 8-3-94

**To calculate sample quantitation limits:
(CAOL * Dilution factor)**

[illegible]

CAOL = Contract Required Quantitation Limit

Action Level Exists

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR 300668

(1/8in)

WATER SAMPLES

**To calculate sample quantitation limits:
(CROL * Dilution factor)**

[illegible]**CRAL = Contract Required Quantitation Limit**

Action Level Exists

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

69900669

Site Name: Westinghouse Sharon

Case #: 22533 Sampling Date(s): 8-2-96 - 8-3-96

WATER SAMPLES (7/58)

To calculate sample quantitation (limits):
(CADL = Dilution Factor)

Sample No.		COA26		COA37	
Dilution Factor		1.0		1.0	
Location		BB-1		BB-2	
COAL	COMPOUND	SAMPLE IS A	SAMPLE IS A	RINSEATE BLANK	RINSEATE BLANK
10	4-Nitroodiphenylamine				
10	4-Bromophenyl phenylether				
10	Hexachlorobenzene				
25	Pentachlorophenol				
10	Phenanthrene				
10	Anthracene				
10	Carbazole				
10	Di-n-butylphthalate				
10	Fluoranthene				
10	Pyrene				
10	Butylbenzylphthalate				
10	3,3'-Dichlorobenzidine				
10	Benzo(a)anthracene				
10	Chrysene				
10	Bis(2-Ethylhexyl)phthalate		3		
10	Di-n-octylphthalate				
10	Benzo(b)fluoranthene				
10	Benzo(k)fluoranthene				
10	Benzo(a)pyrene				
10	Indeno(1,2,3-cd)pyrene				
10	Dibenz(a,h)anthracene				
10	Benzo(g,h,i)perylene				

CRQL = Contract Required Quantitation Limit

Action Level Exists

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR 300670

(b)(7)(D)

To calculate sample quantitation limits:

$$(C_{tol} \times \text{Dilution factor}) / ((100 - X_{moisture}) / 100)$$

Sample No.		COR22	COR23	COR24	COR25	COR27	COR28	COR29	COR30	COR31
Dilution Factor	% Moisture	1.0/2.0	1.0/2.0	1.0	1.0/5.0	1.0	10.0	1.0/10.0	1.0/10.0	1.0
18	4	5	24	19	12	11				
309	\$10	\$11	\$12	\$13	\$14	\$15	\$16	\$17		
CROL	COMPOUND									
330	Hexachlorobutadiene									
330	4-Chloro-3-methylphenol									
330	2-Methylnaphthalene	110	J		200	J		420	560	
330	Hexachlorocyclopentadiene									
330	2,4,6-Trichlorophenol									
800	2,4,5-Trichlorophenol									
330	2-Chloronaphthalene									
800	2-Nitroaniline									
330	Dimethylphthalate									
330	Acenaphthylene							210	200	J
330	2,4-Dinitrotoluene									
800	3-Nitroaniline									
330	Acenaphthene				320	J		380	340	J
800	2,4-Dinitrophenol									R
800	4-Nitrophenol									
330	O benzofuran				280	J		930	740	
330	2,4-Dinitrotoluene								190	J
330	Diethylphthalate									
330	4-Chlorophenyl phenylether									
330	Fluorene	160	J		310	J		420	300	J
800	4-Nitroaniline									
800	4,6-Dinitro-2-methylphenol	100								

CRQL - Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

NR300675

Site Name: Westinghouse Sharon

SOIL SAMPLES
(ug/kg)

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94

To calculate sample quantitation (limits):
(CRQL = Dilution factor / ((100 - Moisture)/100))

Sample No.	CR22	CR23	CR24	CR25	CR27	CR28	CR29	CR30	CR31
Dilution Factor	1.0/2.0	1.0/2.0	1.0	1.0/5.0	1.0	10.0	1.0/10.0	1.0/10.0	1.0
% Moisture	16	6	5	24	19	12	12	11	11
Location	809	810	811	812	813	814	815	816	817
CRQL	COMPOUND								
330	4-Nitroodiphenylamine		UJ						
330	4-Bromophenyl-phenylether		UJ						
330	Hexachlorobenzene		UJ						
800	Pentachlorophenol		UJ	180					
330	Phenanthrene	2300		5300 *	1100	7900	19000 *	16000 *	280
330	Anthracene	320		520	180	1400	900	670	
330	Carbazole	340		500	140		2300	1600	
330	Di-n-butylphthalate			140	770		340	290	450
330	Fluoranthene	3800 *		9300 *	2900	17000	27000 *	24000 *	590
330	Pyrene	2200 *		5700 *	2300	13000	17000 *	16000 *	350
330	Butylbenzylphthalate								
330	5,3'-Dichlorobenzidine								
330	Benzo(a)anthracene	1300		3800 *	890	7700	6600 *	6600 *	240
330	Chrysene	1500		4400 *	970	7900	9100 *	8700 *	270
330	bis(2-Ethylhexyl)phthalate	430		4500 *	620		220		130
330	Di-n-octylphthalate				UJ		UJ	UJ	
330	Benzo(b)fluoranthene	2400		7900 *	1100	9600	8400 *	8600 *	560
330	Benzo(k)fluoranthene	1500		9000	1600	5300	6600 *	6300 *	
330	Benzo(a)pyrene	1400		3400	920	7000	5600 *	5700 *	240
330	Indeno(1,2,3-cd)pyrene	720		800	460	5100	2600	1900	120
330	Benz(a,h)anthracene	380		440	220	1700	920	710	
330	Benzo(g,h,i)perylene	890		1100	520	3300	2800	2000	

CRQL = Contract Required Quantitation Limit

* = Result Reported From Dilution

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR300676

STUDY 1103

(b)(7)(D)

To calculate sample quantitation limits:
 $(CROL * Dilution\ factor) / ((100 - Moisture) / 100)$

[illegible]**CRAL - Contract Required Quantitation Limit**

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR 300677

SOIL SAMPLES
(b)(7)(D)

To calculate sample quantitation limits:
 $(C_{LOL} = \text{dilution factor} / ((100 - \text{moisture})/100))$

[illegible]

CRQL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

AR:300678

SOIL SUPPLIES

(b)(7)(C)

To calculate sample quantitation limits:

$$(CROL * Dilution factor / ((100 - Xmoisture)/100))$$

CROL	COMPOUND	Sample No.		COR32	COR33	COR34	COR35	COR36	COR38	COR39	COR40RE	COR41RE
		Dilution Factor	% Moisture	Location	SAMPLE IS A FIELD DUP. OF	SAMPLE IS B FIELD DUP. OF	SAMPLE IS C FIELD DUP. OF	SAMPLE IS D FIELD DUP. OF	SAMPLE IS E FIELD DUP. OF	SAMPLE IS F FIELD DUP. OF	SAMPLE IS G FIELD DUP. OF	SAMPLE IS H FIELD DUP. OF
330	N-Nitrosodiphenylamine	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	4-Bromophenyl-phenylether	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Hexachlorobenzene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
300	Pentachlorophenol	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Phenanthrene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Anthracene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Carbazole	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Di-n-butylphthalate	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Fluoranthene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Pyrene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Butylbenzylphthalate	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	3,3'-Dichlorobenzidine	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Benzo(a)anthracene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Chrysene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	bis(2-Ethylhexyl)phthalate	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Di-n-octylphthalate	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Benzo(b)fluoranthene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Benzo(k)fluoranthene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Benzo(e)pyrene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Indeno(1,2,3-cd)pyrene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Dibenz(a,h)anthracene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0
330	Benzo(g,h,i)perylene	1.0	17	J	1.0	1.0	1.0/10.0	1.0	1.0	1.0	1.0	1.0

CRCL - Contract Required Quantitation Limit

* = Result Reported From Dilution

♦ = Result Reported from Initial Analysis

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

HR 300679

SOIL SAMPLES
(ug/Kg)

To calculate sample quantitation limits:
 $(C_{CAL} \times \text{Dilution factor}) / ((100 - \% \text{moisture}) / 100)$

Sample No.	COM42	COM43	COM44	COM45	COM46	COM47
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0
% Moisture	13	16	20	12	20	25
Location	S26	S27	S28	S29	S30	S31
CAOL	COMPOUND					
330	Phenol					
330	bis(2-Chloroethyl) ether					
330	2-Chlorophenol					
330	1,3-Dichlorobenzene					
330	1,4-Dichlorobenzene					
330	1,2-Dichlorobenzene					
330	2-Naphthylphenol					
330	2,2'-oxybis(1-chloropropane)					
330	4-Methylphenol					
330	N-Nitroso-di-n-propylamine					
330	Hexachloroethane					
330	Nitrobenzene					
330	Isophorone					
330	2-Nitrophenol					
330	2,4-Dimethylphenol					
330	bis(2-Chloroethoxy)methane					
330	2,4-Dichlorophenol					
330	1,2,4-Trichlorobenzene					
330	Naphthalene	230 +				
330	4-Chloroaniline					

CAOL = Contract Required Quantitation Limit

† = Result Reported From Reanalysis

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

AR 300680

Site Name: Westinghouse Sharon

SOIL SAMPLES
(ug/Kg)

Case #: Z2533 Sampling Date(s): 8-2-94 - 8-3-94

To calculate sample quantitation limits:
(CROL * Dilution factor / ((100 - %moisture)/100)

Sample No.	COR42	COR43	COR44	COR45	COR46	COR47
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0
% Moisture	13	16	20	12	20	25
Location	326	327	328	329	330	331
CROL	COMPOUND					
330	Hexachlorobutadiene					
330	4-Chloro-3-methylphenol					
330	2-Methylnaphthalene	310 + J				
330	Hexachlorocyclopentadiene					
330	2,4,6-Trichlorophenol					
800	2,4,5-Trichlorophenol					
330	2-Chloronaphthalene					
800	2-Nitroaniline					
330	Dimethylphthalate					
330	Acenaphthylene					
330	2,6-Dinitrotoluene					
800	3-Nitroaniline					
330	Acenaphthene					
800	2,4-Dinitrophenol	100 + J				
800	4-Nitrophenol					
330	Dibenzofuran					
330	2,4-Dinitrotoluene					
330	Diethylphthalate					
330	4-Chlorophenyl-phenylether					
330	Fluorene					
800	4-Nitroaniline					
800	4,6-Dinitro-2-methylphenol					

CROL = Contract Required Quantitation Limit

+ = Result Reported From Reanalysis

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR3000684

Site Name: Westinghouse Shoran

SOIL SAMPLES

(ug/Kg)

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94

To calculate sample quantitation limits:
(CROL = Dilution factor / ((100 - Moisture)/100))

Sample No.	COR42	COR43	COR44	COR45	COR46	COR47
Dilution factor	1.0	1.0	1.0	1.0	1.0	1.0
% Moisture	13	16	20	12	20	25
Location	S26	S27	S28	S29	S30	S31
CROL	COMPOUND					
330	N-Nitrosodiphenylamine					
330	4-Bromophenyl-phenylether					
330	Hexachlorobenzene					
800	Pentachlorophenol					
330	Phenanthrene	640	670			
330	Anthracene	130	J			
330	Carbazole	130	J	120		99
330	Di-n-butylphthalate	1900	1200	820	2400	2200
330	Fluoranthene	1900	270	J	630	190
330	Pyrene	940	J	1300	540	J
330	Butylbenzylphthalate					J
330	3,3'-Dichlorobenzidine					J
330	Benzo(a)anthracene	940	J	930	470	J
330	Chrysene	780	J	950	400	J
330	bis(2-Ethylhexyl)phthalate	130 +	J	130	160	120
330	Di-n-octylphthalate					J
330	Benzo(b)fluoranthene	890	J	1100	550	170
330	Benzo(k)fluoranthene	1100	J	1200	550	J
330	Benzo(a)pyrene	650	J	1000	480	J
330	Indeno(1,2,3-cd)pyrene	1200	J	690	430	J
330	Dibenz(a,h)anthracene	1100	J	280	130	J
330	Benzo(g,h,i)perylene	1300	J	760	660	J

CROL = Contract Required Quantitation Limit

+ = Result Reported From Reanalysis

SEE NARRATIVE FOR CODE DEFINITIONS
revised 07/90

AR300682

WATER SAMPLES

To calculate sample quantitation limits:
(CNDL = Dilution Factor)

Sample No.		CON26		COR37	
Dilution Factor	Location	1.0	1.0		
		R8-1	R8-2		
COROL	COMPOUND	SAMPLE IS A RINSEATE BLANK. RINSEATE BLANK.	SAMPLE IS A		
	alpha-BHC	[W]	[W]		
	beta-BHC	[W]	[W]		
	delta-BHC	[W]	[W]		
	*gamma-BHC (Lindane)	[W]	[W]		
	*Heptachlor	[W]	[W]		
	Aldrin	[W]	[W]		
	Neptechlor Epoxide	[W]	[W]		
	Endosulfan I	[W]	0.016 J		
	Bieldrin	[W]	[W]		
	4,4'-DDE	[W]	0.0064 J		
	*Endrin	[W]	[W]		
	Endosulfan II	[W]	[W]		
	4,4'-DDD	[W]	[W]		
	Endosulfan Sulfate	[W]	[W]		
	4,4'-DDT	[W]	[W]		
	*Methoxychlor	[W]	[W]		
	Endrin Ketone	[W]	[W]		
	Endrin Aldehyde	[W]	[W]		
	*alpha-Chlordane	[W]	0.018 J		
	*gamma-Chlordane	[W]	[W]		
	*Toxaphene	[W]	[W]		
	*Aroclor-1016	[W]	[W]		
	*Aroclor-1221	[W]	[W]		
	*Aroclor-1232	[W]	[W]		
	*Aroclor-1242	[W]	[W]		
	*Aroclor-1248	[W]	[W]		
	*Aroclor-1254	[W]	[W]		
	*Aroclor-1260	[W]	[W]		

CRQL = Contract Required Quantitation Limit

44-38861-103

revised 07/90

Site Name: Westinghouse Sharon

SOIL SAMPLES

(ug/Kg)

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94

To calculate sample quantitation limits:
(CROL = dilution factor / ((100 - %moisture)/100))

Sample No.	COR13	COR14	COR15	COR16	COR17	COR18	COR19	COR20	COR21
Dilution Factor	1.0/10.0	1.0	1.0	2.0/20.0	1.0	1.0	1.0	1.0	1.0
% Moisture	7	21	10	23	8	19	15	21	12
Location	S01	S02	S03	S04	S05	S06	S06 DUPE	S07	S08
COMPOUND						SAMPLE IS A FIELD DUP. OF COR18			
1.7	alpha-BHC								
1.7	beta-BHC								
1.7	delta-BHC								
1.7	gamma-BHC (Lindane)								
1.7	Heptachlor								
1.7	Aldrin								
1.7	Heptachlor Epoxide								
1.7	Endosulfan I								
3.3	Dieldrin	300*			8.3				
3.3	4,4'-DDE	7.6	26	82	7.3	7.3	11	12	2.1
3.3	Endrin		15	9.5		18	20		5.0
3.3	Endosulfan II								
3.3	4,4'-DDD		30						
3.3	Endosulfan Sulfate								
3.3	4,4'-DDT	30	21		170	26	42	25	7.4
17	Methoxychlor								
3.3	Endrin Ketone								
3.3	Endrin Aldehyde			11		27	46	24	7.0
1.7	alpha-Chlordane				4.7	51	60	2.4	24
1.7	gamma-Chlordane				5.0	38	46	1.4	16
170	Toxaphene								
33	Aroclor-1016								
67	Aroclor-1221								
33	Aroclor-1232								
33	Aroclor-1242								
33	Aroclor-1248								
33	Aroclor-1254								
33	Aroclor-1260	410	1300	440	980	840	1400	630	220
CROL	Contract Required Quantitation Limit								

* = Results taken from diluted analyses

AR300684

DATA SUMMARY FORM: PESTICIDES AND PCBS

Page 28 of 30

Site Name: Westinghouse Shoron

SOIL SAMPLES

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94

To calculate sample quantitation limits:
(CROL = Dilution factor / ((100 - Moisture)/100))

Sample No.	COR22	COR23	COR24	COR25	COR27	COR28	COR29	COR30	COR31
Dilution Factor	1.0	1.0	20.0	1.0	10.0	1.0	2.0	2.0	1.0
% Moisture	18	4	5	24	19	12	12	11	11
Location	S10	S10	S11	S12	S13	S14	S15	S16	S17
CROL	COMPOUND								
1.7	alpha-BHC								
1.7	beta-BHC								
1.7	delta-BHC								
1.7	gamma-BHC (Lindane)								
1.7	Heptachlor								
1.7	Aldrin								
1.7	Heptachlor Epoxide								
1.7	Endosulfan I								
3.3	Dieldrin	6.1					41		
3.3	4,4'-DDE	15				15			
3.3	Endrin	8.4				12			
3.3	Endosulfan II							70	1.2
3.3	4,4'-DDD								
3.3	Endosulfan Sulfate								
3.3	4,4'-DDT	56							
17	Methoxychlor						62	37	2.7
3.3	Endrin Ketone								
3.3	Endrin Aldehyde	27							
1.7	alpha-Chlordane	9.4							
1.7	gamma-Chlordane	8.3							
170	Toxaphene								
33	Aroclor-1016								
67	Aroclor-1221								
33	Aroclor-1232								
33	Aroclor-1242								
33	Aroclor-1248								
33	Aroclor-1254								
33	Aroclor-1260	760	870	550	1200	800		1800	76

CROL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

AR3000685

revised 07/90

DATA SUMMARY FORM: PESTICIDES AND PCBs

Page 29 of 30

Site Name: Westinghouse Shoron

SOIL SAMPLES

(ug/Kg)

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94

To calculate sample quantitation limits:
(CAL = Dilution factor / ((100 - Moisture)/100))

Sample No.	COR32	COR33	COR34	COR35	COR36	COR38	COR39	COR40	COR41
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0	1.0
% Moisture	17	15	15	28	18	15	14	11	15
Location	S14	S19	S19 DUPE	S20	S21	S22	S23	S24	S25
	SAMPLE 18 A	SAMPLE 18 A	SAMPLE 18 A						
	FIELD DUP. OF	FIELD DUP. OF	FIELD DUP. OF						
	COR34	COR33	COR33						
CONC	CONC	CONC	CONC	CONC	CONC	CONC	CONC	CONC	CONC
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
alpha-BHC	alpha-BHC	alpha-BHC	alpha-BHC	alpha-BHC	alpha-BHC	alpha-BHC	alpha-BHC	alpha-BHC	alpha-BHC
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
beta-BHC	beta-BHC	beta-BHC	beta-BHC	beta-BHC	beta-BHC	beta-BHC	beta-BHC	beta-BHC	beta-BHC
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
delta-BHC	delta-BHC	delta-BHC	delta-BHC	delta-BHC	delta-BHC	delta-BHC	delta-BHC	delta-BHC	delta-BHC
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
gamma-BHC (Lindane)	gamma-BHC (Lindane)	gamma-BHC (Lindane)	gamma-BHC (Lindane)	gamma-BHC (Lindane)	gamma-BHC (Lindane)	gamma-BHC (Lindane)	gamma-BHC (Lindane)	gamma-BHC (Lindane)	gamma-BHC (Lindane)
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
Heptachlor	Heptachlor	Heptachlor	Heptachlor	Heptachlor	Heptachlor	Heptachlor	Heptachlor	Heptachlor	Heptachlor
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
Aldrin	Aldrin	Aldrin	Aldrin	Aldrin	Aldrin	Aldrin	Aldrin	Aldrin	Aldrin
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
Heptachlor Epoxide	Heptachlor Epoxide	Heptachlor Epoxide	Heptachlor Epoxide	Heptachlor Epoxide	Heptachlor Epoxide	Heptachlor Epoxide	Heptachlor Epoxide	Heptachlor Epoxide	Heptachlor Epoxide
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
Endosulfan I	Endosulfan I	Endosulfan I	Endosulfan I	Endosulfan I	Endosulfan I	Endosulfan I	Endosulfan I	Endosulfan I	Endosulfan I
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
Bifenthrin	Bifenthrin	Bifenthrin	Bifenthrin	Bifenthrin	Bifenthrin	Bifenthrin	Bifenthrin	Bifenthrin	Bifenthrin
3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3
4,4'-DDE	4,4'-DDE	4,4'-DDE	4,4'-DDE	4,4'-DDE	4,4'-DDE	4,4'-DDE	4,4'-DDE	4,4'-DDE	4,4'-DDE
3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3
Endrin	Endrin	Endrin	Endrin	Endrin	Endrin	Endrin	Endrin	Endrin	Endrin
3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3
Endosulfan II	Endosulfan II	Endosulfan II	Endosulfan II	Endosulfan II	Endosulfan II	Endosulfan II	Endosulfan II	Endosulfan II	Endosulfan II
3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3
4,4'-DDD	4,4'-DDD	4,4'-DDD	4,4'-DDD	4,4'-DDD	4,4'-DDD	4,4'-DDD	4,4'-DDD	4,4'-DDD	4,4'-DDD
3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3
Endosulfan Sulfate	Endosulfan Sulfate	Endosulfan Sulfate	Endosulfan Sulfate	Endosulfan Sulfate	Endosulfan Sulfate	Endosulfan Sulfate	Endosulfan Sulfate	Endosulfan Sulfate	Endosulfan Sulfate
3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3
4,4'-DDT	4,4'-DDT	4,4'-DDT	4,4'-DDT	4,4'-DDT	4,4'-DDT	4,4'-DDT	4,4'-DDT	4,4'-DDT	4,4'-DDT
17	17	17	17	17	17	17	17	17	17
Metobuthylchlor	Metobuthylchlor	Metobuthylchlor	Metobuthylchlor	Metobuthylchlor	Metobuthylchlor	Metobuthylchlor	Metobuthylchlor	Metobuthylchlor	Metobuthylchlor
3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3
Endrin Ketone	Endrin Ketone	Endrin Ketone	Endrin Ketone	Endrin Ketone	Endrin Ketone	Endrin Ketone	Endrin Ketone	Endrin Ketone	Endrin Ketone
3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3	3.3
Endrin Aldehyde	Endrin Aldehyde	Endrin Aldehyde	Endrin Aldehyde	Endrin Aldehyde	Endrin Aldehyde	Endrin Aldehyde	Endrin Aldehyde	Endrin Aldehyde	Endrin Aldehyde
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
alpha-Chlordane	alpha-Chlordane	alpha-Chlordane	alpha-Chlordane	alpha-Chlordane	alpha-Chlordane	alpha-Chlordane	alpha-Chlordane	alpha-Chlordane	alpha-Chlordane
1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7	1.7
gamma-Chlordane	gamma-Chlordane	gamma-Chlordane	gamma-Chlordane	gamma-Chlordane	gamma-Chlordane	gamma-Chlordane	gamma-Chlordane	gamma-Chlordane	gamma-Chlordane
170	170	170	170	170	170	170	170	170	170
Tenaphene	Tenaphene	Tenaphene	Tenaphene	Tenaphene	Tenaphene	Tenaphene	Tenaphene	Tenaphene	Tenaphene
33	33	33	33	33	33	33	33	33	33
Aroclor-1016	Aroclor-1016	Aroclor-1016	Aroclor-1016	Aroclor-1016	Aroclor-1016	Aroclor-1016	Aroclor-1016	Aroclor-1016	Aroclor-1016
67	67	67	67	67	67	67	67	67	67
Aroclor-1221	Aroclor-1221	Aroclor-1221	Aroclor-1221	Aroclor-1221	Aroclor-1221	Aroclor-1221	Aroclor-1221	Aroclor-1221	Aroclor-1221
33	33	33	33	33	33	33	33	33	33
Aroclor-1232	Aroclor-1232	Aroclor-1232	Aroclor-1232	Aroclor-1232	Aroclor-1232	Aroclor-1232	Aroclor-1232	Aroclor-1232	Aroclor-1232
33	33	33	33	33	33	33	33	33	33
Aroclor-1242	Aroclor-1242	Aroclor-1242	Aroclor-1242	Aroclor-1242	Aroclor-1242	Aroclor-1242	Aroclor-1242	Aroclor-1242	Aroclor-1242
33	33	33	33	33	33	33	33	33	33
Aroclor-1248	Aroclor-1248	Aroclor-1248	Aroclor-1248	Aroclor-1248	Aroclor-1248	Aroclor-1248	Aroclor-1248	Aroclor-1248	Aroclor-1248
33	33	33	33	33	33	33	33	33	33
Aroclor-1254	Aroclor-1254	Aroclor-1254	Aroclor-1254	Aroclor-1254	Aroclor-1254	Aroclor-1254	Aroclor-1254	Aroclor-1254	Aroclor-1254
33	33	33	33	33	33	33	33	33	33
Aroclor-1260	Aroclor-1260	Aroclor-1260	Aroclor-1260	Aroclor-1260	Aroclor-1260	Aroclor-1260	Aroclor-1260	Aroclor-1260	Aroclor-1260
250	250	250	250	250	250	250	250	250	250

CAL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

AR300686

revised 07/90

Site Name: Westinghouse Sharon

SOIL SAMPLES

Case #: 22533 Sampling Date(s): 8-2-94 - 8-3-94 To calculate sample quantitation limits:
(CROL = Dilution factor / ((100 - Moisture)/100))

Sample No.	COR42	COR43	COR44	COR45	COR46	COR47
Dilution Factor	1.0	1.0	1.0	1.0	1.0	1.0
% Moisture	13	16	20	12	20	25
Location	S26	S27	S28	S29	S30	S31
CROL	COMPOUND					
1.7	alpha-BHC	UL	UL	UL	UL	UL
1.7	Beta-BHC	UL	UL	UL	UL	UL
1.7	delta-BHC	UL	UL	UL	UL	UL
1.7	gamma-BHC (Lindane)	UL	UL	UL	UL	UL
1.7	Heptachlor	UL	UL	UL	UL	UL
1.7	Aldrin	UL	UL	UL	UL	UL
1.7	Heptachlor Epoxide	0.87	0.66	1.1	UL	UL
1.7	Endosulfan I	3.1	17	1.9	3.3	1.2
3.3	Dieldrin	UL	UL	UL	UL	UL
3.3	4,4'-DDE	UL	7.1	UL	1.1	0.85
3.3	Endrin	UL	UL	UL	UL	UL
3.3	Endosulfan II	UL	UL	UL	UL	UL
3.3	4,4'-DDD	UL	UL	UL	UL	UL
3.3	Endosulfan Sulfate	UL	UL	UL	UL	UL
3.3	4,4'-DDT	1.3	5.3	UL	0.67	UL
17	Methoxychlor	UL	UL	UL	UL	UL
3.3	Endrin Ketone	UL	UL	UL	UL	UL
3.3	Endrin Aldehyde	1.7	6.6	2.4	UL	UL
1.7	alpha-Chlordane	2.7	5.4	1.2	2.4	UL
1.7	gamma-Chlordane	UL	UL	UL	UL	UL
170	Toxaphene	UL	UL	UL	UL	UL
33	Aroclor-1016	UL	UL	UL	UL	UL
67	Aroclor-1221	UL	UL	UL	UL	UL
33	Aroclor-1232	UL	UL	UL	UL	UL
33	Aroclor-1242	UL	UL	UL	UL	UL
33	Aroclor-1248	UL	UL	UL	UL	UL
33	Aroclor-1254	UL	UL	UL	UL	UL
33	Aroclor-1260	UL	100	UL	UL	17

CROL = Contract Required Quantitation Limit

SEE NARRATIVE FOR CODE DEFINITIONS

revised 07/90

AR300687



United States Environmental Protection Agency
Contract Laboratory Program

Organic Traffic Report & Chain of Custody Record

(For Organic CLP Analysis)

Case No.

SAS No.
(if applicable)

22533

1. Project Code:
5303

2. Region No. Sampling Co.
III WESTON

4. Date Shipped Carrier
8/4/94 Federal Express

Regional Information

3. Purpose:
SF
ST
FED

5. Ship To
Envirosystems, Inc.
9200 Rumsy Rd, Suite # B102
Columbia, MD 21045
ATTN: Mohan Khare

6. Matrix
(Enter in Column A)

7. Preservative
(Enter in Column D)

Non-Superfund Program

8. Matrix
(Enter in Column A)

9. Other (Specify in Column A)

Site Name

Westinghouse Sharon

City, State

Sharon, PA

10. Matrix
(Enter in Column A)

11. Other (Specify in Column A)

12. Other (Specify in Column A)

13. Other (Specify in Column A)

14. Other (Specify in Column A)

15. Other (Specify in Column A)

16. Other (Specify in Column A)

17. Other (Specify in Column A)

18. Other (Specify in Column A)

19. Other (Specify in Column A)

20. Other (Specify in Column A)

21. Other (Specify in Column A)

22. Other (Specify in Column A)

23. Other (Specify in Column A)

24. Other (Specify in Column A)

25. Other (Specify in Column A)

26. Other (Specify in Column A)

27. Other (Specify in Column A)

28. Other (Specify in Column A)

29. Other (Specify in Column A)

30. Other (Specify in Column A)

31. Other (Specify in Column A)

32. Other (Specify in Column A)

33. Other (Specify in Column A)

34. Other (Specify in Column A)

35. Other (Specify in Column A)

36. Other (Specify in Column A)

37. Other (Specify in Column A)

38. Other (Specify in Column A)

39. Other (Specify in Column A)

40. Other (Specify in Column A)

41. Other (Specify in Column A)

42. Other (Specify in Column A)

43. Other (Specify in Column A)

44. Other (Specify in Column A)

45. Other (Specify in Column A)

46. Other (Specify in Column A)

47. Other (Specify in Column A)

48. Other (Specify in Column A)

DISTRICT:

Blue - Region Copy

White - Lab Copy for Return to Region

Pink - SMO Copy

Yellow - Lab Copy for Return to

EPA Form 910-3

SEE REVERSE FOR ADDITIONAL STANDARD

*SEE REVERSE FOR PURPOSE CODE DEFINITION

CTIONS

360850

AR300688



United States Environmental Protection Agency
Contract Laboratory Program

Organic Traffic Report & Chain of Custody Record

(For Organic CLP Analysis)

Case No.

22533

SAS No.
(if applicable)

6. Matrix
(Enter in Column A)

- 1. Surface Water
- 2. Ground Water
- 3. Leachate
- 4. Field QC
- 5. Soil/Sediment
- 6. Oil (High only)
- 7. Waste (High only)
- 8. Other (Specify in Column A)

7. Preservative
(Enter in Column D)

- 1. HCl
- 2. HNO₃
- 3. NaHSO₄
- 4. H₂SO₄
- 5. Ice only
- 6. Other (Specify in Column D)
- N. Not preserved

4. Date Shipped (Carrier)

8/4/94 Federal Express

Airbill Number

202 868 0953

5. Ship To

Enviro Systems Inc.
9200 Rumsby Rd, Suite #B102
Columbia, MD 21045
ATTN: Mohan Khare

2. Region No. (Sampling Co.)

WESTON

Sampler (Name)

THE CARTER

Sampler Signature

Lead ☒ CLEM ☐ PA ☐ REM ☐ RI ☐ SI ☐ ESI ☐ NPLD

1. Project Code

5303

Account Code

Site Name

Westinghouse - Sharon

City, State
Sharon, PA

CLP Sample Numbers (from labels)

Matrix (from Box 6) Other

Conc.: Low Med High

Sample Type: Grab (from Box 7) Other

Preservative (from Box 7) Other

Regional Specific Tracking Number or Tag Numbers

Station Location Identifier

Mo/Day/Year/Time Sample Collection

Corresponding CLP Inorganic Sample No.

Sampler Initials

Field QC Qualifier
B = Blank S = Spike
D = Duplicate
R = Retest
HE = Perform Eval
- = No QC Sample

Shipment for Case Complete? (Y/N)

Page 4 of 4

Sample(s) to be Used for Laboratory QC

Additional Samples Signatures

Chain of Custody Seal Number(s)

CHAIN OF CUSTODY RECORD

Relinquished by: (Signature) C. V. Adell	Date / Time 8/4/94 1600	Received by: (Signature)	Date / Time
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time
Relinquished by: (Signature)	Date / Time	Received by: (Signature)	Date / Time

DISTRIBUTION: Blue - Region Copy White - Lab Copy for Return to Region Yellow - Lab Copy for Return to SMO

EPA Form 8110-2

SEE REVERSE FOR ADDITIONAL STANDARD INSTRUCTIONS
SEE REVERSE FOR PURPOSE CODE DEFINITIONS

AR300691

Attachment B

**Dioxin/Dibenzo furan Data Validation, Analytical Summary,
Chain-of-Custody**

AR300692



5 Underwood Court, Delran, New Jersey 08075-1229
609-461-4003 • 215-238-0338 • Fax 609-461-4916

TECHNICAL ASSISTANCE TEAM FOR EMERGENCY RESPONSE REMOVAL AND PREVENTION
EPA CONTRACT 68-WO-0036

MEMORANDUM

TO: Gerald Heston, OSC, EPA Region III
Western Removal Section

FROM: Marian Murphy, TAT Region III ~~AW~~ TDD# 9205-25B
PCS# 5497

SUBJECT: Westinghouse Sharon Samples Analytical Review

DATE: September 15, 1994

This report covers the general review of the data package submitted by Weston Lionville Analytical Laboratory, for one (1) rinsate blank sample and fifteen (15) soil samples collected at the Westinghouse Sharon Site on August 2-3, 1994. The samples were received at Weston Lionville Analytical Laboratory, in Lionville, PA on August 6, 1994. The analysis requested was isomer specific polychlorinated dibenzo dioxins and polychlorinated dibenzo furans and to determine the toxicity equivalents as 2,3,7,8 tetra chloro dibenzo dioxin (2,3,7,8-TCDD).

ANALYTICAL METHODOLOGY

The samples were analyzed for isomer specific polychlorinated dibenzo dioxin and polychlorinated dibenzo furans following EPA Method 8280.

- Signed chain-of-custody records were received.
- The hold times were met. The initial and continuing calibration data met ion ratio criteria, percent relative standard deviation criteria and percent difference criteria, respectively. The method blanks were free of contamination. The surrogate spike recoveries met criteria. The MS/MSD recoveries and the RPD values met criteria. All identified isomers met identification criteria. Attached is a table of 2,3,7,8-TCDD toxicity equivalent factors for all samples which positive values for any isomer or homolog were reported.

CONCLUSION

Accept all data as presented.

Attachment: 2,3,7,8-TCDD Toxicity Equivalents Factors

Roy F. Weston, Inc.
MAJOR PROGRAMS DIVISION
In Association with Foster Wheeler USA Corporation, Resource Applications, Inc., C.C. Johnson & Malhotra, P.C.,
R.E. Sarriera Associates, and GRB Environmental Services, Inc.

AR300693

2.3.7.8-TCDD TOXICITY EQUIVALENTS FOR THE WESTINGHOUSE SHARON SITE

UNITS EXPRESSED AS (ng/g) OR PARTS PER BILLION

ISOMER/HOMOLOG	S01	S04	S06	S07	S08	S16	S16	S16	S24	S27	S40	S51
	NG/G	NG/G	NG/G	NG/G	NG/G	NG/G	NG/G	NG/G	NG/G	NG/G	NG/G	NG/G
1,2,3,6,7,8-HxCDD			0.0008									
TOTAL HxCDD		0.0024	0.0028	0.0016								
1,2,3,4,6,7,8-HxCDF	0.0001	0.0009	0.0001	0.0002	0.0001							
TOTAL HxCDF	0.0001	0.0009	0.0001	0.0002	0.0001							
OCDD	0.00005	0.0005	0.0005	0.0007	0.0005	0.0004	0.0004	0.0004	0.0008	0.0009	0.0006	
2,3,7,8-TCDF		0.003	0.003	0.003								
TOTAL TCDF		0.0001	0.00014	0.00011								
TOTAL PCDF	0.005	0.018	0.022	0.008								
1,2,3,4,7,8-HxCDF	0.0001	0.0003	0.0003	0.0002								
1,2,3,6,7,8-HxCDF		0.0001	0.0001									
2,3,4,6,7,8-HxCDF		0.0002	0.0002									
TOTAL HxCDF	0.0009	0.0018	0.0021	0.0011								
1,2,3,4,6,7,8-HPCDF	0.00004	0.00006	0.00006	0.00006			0.002					
TOTAL HPCDF	0.00008	0.00006	0.00006	0.00005								
OCDF	0.0001	0.00007	0.0001	0.0001			0.0001					
TOTAL 2,3,7,8-TEQ	0.00547	0.02876	0.03235	0.01322	0.0035	0.0004	0.0026	0.0004	0.0006	0.0009	0.0006	0.0006

AR300694

PROJ. NO. 5303 PROJECT NAME Sharon Westinghouse
 SAMPLERS: (Signature) Joseph A. Carter

STA. NO.	DATE	TIME	NO. OF CONTAINERS	STATION LOCATION	REMARKS
S01	8/2/94	1445	1	S01	
S04	8/2/94	1810	2	Blender Rinse (amber)	
S06	8/2/94	1420	1	S04	
S07	8/2/94	1510	1	S06	
S09	8/2/94	1535	1	S07	
S15	8/2/94	1630	1	S09	
S16	8/3/94	1015	1	S15	
S17	8/3/94	1030	1	S16	
S20	8/3/94	1055	1	S17	
S24	8/3/94	1200	1	S20	
S27	8/3/94	1535	1	S24	
S28	8/3/94	1645	1	S27	
S29	8/3/94	1735	1	S28	
S30	8/3/94	1800	1	S29	
S30	8/3/94	1845	1	S30	

Relinquished by: (Signature) Joseph A. Carter Date / Time Received by: (Signature) Date / Time

Relinquished by: (Signature) Date / Time Received by: (Signature) Date / Time

Relinquished by: (Signature) Date / Time Received by: (Signature) Date / Time

Remarks

HR 300695

Distribution: Original Accompanies Shipment; Copy to Coordinator Field